# Final Third Quarter 2014 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells

Red Hill Bulk Fuel Storage Facility
Joint Base Pearl Harbor-Hickam, Oahu, Hawaii

DOH Facility ID: 9-102271

DOH Release ID: 990051, 010011, 020028, and 140010

September 2014

Department of the Navy Naval Facilities Engineering Command, Hawaii 400 Marshall Road JBPHH HI 96860-3139



Contract Number N62742-12-D-1853, CTO 0002



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Prepared for:



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Prepared by:

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Prepared under:

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# FINAL THIRD QUARTER 2014 - QUARTERLY GROUNDWATER MONITORING REPORT INSIDE TUNNEL WELLS RED HILL BULK FUEL STORAGE FACILITY

Long-Term Groundwater and Soil Vapor Monitoring Red Hill Bulk Fuel Storage Facility Joint-Base Pearl Harbor-Hickam, Oahu, Hawaii

### **Prepared for:**

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Contract Number: N62742-12-D-1853 Contract Task Order: 0002

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### **ACRONYMS AND ABBREVIATIONS**

ACRONYMS/

ABBREVIATIONS DEFINITION/MEANING

% percent

bgs below ground surface

COPC Contaminant of Potential Concern

DLNR State of Hawaii Department of Land and Natural Resources

DOH State of Hawaii Department of Health

DON Department of the Navy
EAL Environmental Action Level
EPA Environmental Protection Agency
ESI Environmental Science International

F-76 Marine Diesel Fuel ID Identification

JBPHH Joint Base Pearl Harbor-Hickam

JP-5 Jet Fuel Propellant-5
JP-8 Jet Fuel Propellant-8
LCS Laboratory Control Sample

LCSD Laboratory Control Sample Duplicate

LOD Limit of Detection
LOQ Limit of Quantitation

µg/L micrograms per liter

MS Matrix Spike

MSD Matrix Spike Duplicate

NAVFAC Naval Facilities Engineering Command

NAVSUP FLC Naval Supply Systems Command Fleet Logistics Center

PAH Polycyclic Aromatic Hydrocarbons

PARCCS Precision, Accuracy, Representativeness, Completeness, Comparability,

and Sensitivity

pH hydrogen activity

RHSF Red Hill Bulk Fuel Storage Facility

RPD Relative Percent Difference
SAP Sampling and Analysis Plan
SSRBL Site-Specific Risk-Based Level
TEC The Environmental Company, Inc.

TPH-d Total Petroleum Hydrocarbons as diesel TPH-g Total Petroleum Hydrocarbons as gasoline

U.S. United States of America
UST Underground Storage Tank
VOC Volatile Organic Compound

WP Work Plan

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#### **EXECUTIVE SUMMARY**

This quarterly groundwater monitoring report presents the results of the third quarter 2014 groundwater sampling conducted on July 21 and 22, 2014, at the Red Hill Bulk Fuel Storage Facility [RHSF], Joint Base Pearl Harbor-Hickam [JBPHH], Hawaii. The RHSF is located in Halawa Heights on the Island of Oahu. There are 18 active and 2 inactive Underground Storage Tanks [USTs] located at the RHSF. The State of Hawaii Department of Health [DOH] Facility Identification [ID] number is 9-102271. The DOH Release ID numbers are 990051, 010011, 020028, and 140010.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring at the RHSF and concurrent with release response activities initiated at Tank 5 in January, under Naval Facilities Engineering Command [NAVFAC] Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved 2012 Work Plan [WP]/Sampling and Analysis Plan [SAP] prepared by Environmental Science International, Inc. [ESI].

On July 21 and 22, 2014, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01). A primary and duplicate sample were collected from well RHMW02. A summary of the analytical results is provided below.

- RHMW01 Total Petroleum Hydrocarbons as diesel fuel [TPH-d] (67 micrograms per liter [µg/L]) was the only analyte detected. The TPH-d concentration detected did not exceed the DOH Environmental Action Levels [EALs] or the site-specific risk-based level [SSRBL]. TPH-d concentrations in this well have shown an overall decreasing trend from a high of 1,500 µg/L in February 2005.
- RHMW02 TPH-d (1,200 and 1,300 μg/L), Total Petroleum Hydrocarbons as gasoline [TPH-g] (48 and 49  $\mu$ g/L), xylenes (0.36 and 0.33  $\mu$ g/L), acenaphthene (0.52 and 0.50  $\mu$ g/L), fluorene (0.24 and 0.23 μg/L), 1-methylnaphthalene (25 and 26 μg/L), 2-methylnaphthalene (20 and 22 μg/L), and naphthalene (71 and 76 μg/L) were detected. TPH-d was detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination, but below half the SSRBL. The polycyclic aromatic hydrocarbons [PAHs] 1-methylnaphthalene and naphthalene were detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination. 2-Methylnaphthalene was detected at concentrations above the DOH EAL for gross contamination. TPH-d concentrations have generally been decreasing in this well since January 15, 2014; however. the concentrations of 1-methylnaphthalene, 2-methylnaphthalene, naphthalene increased to their highest levels since January 2013.
- RHMW03 TPH-d (37 μg/L) was the only analyte detected. None of the chemical constituents analyzed for were detected at concentrations above the DOH EALs.

- RHMW05 None of the chemical constituents analyzed for were detected.
- RHMW2254-01 None of the chemical constituents analyzed for were detected.

Between January and July 2014, the sampling frequency of the wells and monitoring point inside Red Hill increased, due to response activities related to a reported release at Tank 5. Since the previous quarterly sampling event in April 2014, three wells (RHMW01, RHMW02, and RHMW05) and one sampling point (RHMW2254-01) have each been sampled three times, including this quarterly event. During this period only TPH-d in RHMW01 and TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 were detected at concentrations exceeding the DOH EALs. However, the TPH-d concentration detected in RHMW01 only exceeded the DOH EAL during the April 7, 2014 release response groundwater monitoring event. TPH-d concentrations were less than half the SSRBL. Groundwater contaminant concentrations for this period in the other three wells (RHMW03, RHMW05, and RHMW2254-01) remained at low concentrations and did not change significantly, or were not detected.

Concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 show an increasing trend since March 5, 2014, but remain below the historic maximums detected in the well. All other analytical results were generally consistent with historical data.

Based on the groundwater monitoring results and the reported release at Tank 5 in January 2014, continued groundwater monitoring at the wells inside the RHSF tunnel is recommended. The next quarterly event is tentatively scheduled for October 2014.

#### **SECTION 1 – INTRODUCTION**

This quarterly groundwater monitoring report presents the results of the third quarter 2014 groundwater sampling conducted on July 21 and 22, 2014, at the RHSF, JBPHH. The RHSF is located in Halawa Heights on the Island of Oahu. The purpose of the sampling is to (1) assess the condition of groundwater beneath the RHSF with respect to chemical constituents associated with jet fuel propellant and marine diesel fuel, and (2) to ensure the Navy remains in compliance with DOH UST release response requirements as described in Hawaii Administrative Rules Chapter 11-281 Subchapter 7, *Release Response Action* (DOH, 2013). The DOH Facility ID number for the RHSF is 9-102271. The DOH Release ID numbers are 990051, 010011, 020028, and 140010.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI (ESI, 2012).

#### 1.1 SITE DESCRIPTION

The RHSF is located on federal government land (zoned F1- Military and Federal), located in Halawa Heights, approximately 2.5 miles northeast of Pearl Harbor (Figure 1). It is located on a low ridge on the western edge of the Koolau Mountain Range that divides Halawa Valley from Moanalua Valley. The RHSF is bordered on the north by Halawa Correctional Facility and private businesses, on the west by the United States of America [U.S.] Coast Guard reservation, on the south by residential neighborhoods, and on the east by Moanalua Valley. A quarry is located less than a quarter mile away to the northwest. The RHSF occupies 144 acres of land and the majority of the site is at an elevation of approximately 200 to 500 feet above mean sea level.

The RHSF contains 18 active and 2 inactive USTs that are operated by Naval Supply Systems Command Fleet Logistics Center [NAVSUP FLC] Pearl Harbor (formerly Fleet and Industrial Supply Center). Each UST has a capacity of approximately 12.5 million gallons. The RHSF is located approximately 100 feet above the basal aquifer. The USTs contain Jet Fuel Propellant-5 [JP-5], Jet Fuel Propellant-8 [JP-8], and Marine Diesel Fuel [F-76]. The current status of the USTs are summarized in Table 1.1.

Four groundwater monitoring wells (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01) are located within the RHSF lower access tunnel (Figure 2). Three groundwater monitoring wells (wells HDMW2253-03, OWDFMW01, and RHMW04) are located outside of the RHSF tunnel system. Monitoring data for the three wells located outside the tunnel are included in a separate report.

As noted, monitoring wells RHMW01, RHMW02, RHMW03, and RHMW05 are located inside the underground tunnels. Sampling point RHMW2254-01 is located inside the infiltration gallery of the Department of the Navy [DON] drinking water supply Well 2254-01. The DON Well

2254-01 is located approximately 2,400 feet downgradient of the USTs and provides potable water to the JBPHH Water System, which serves approximately 65,200 military customers. NAVFAC Public Works Department operates the infiltration gallery and DON Well 2254-01.

TABLE 1.1
Current Status of the USTs
Red Hill Bulk Fuel Storage Facility
July 2014 Quarterly Monitoring Report

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Tank Identification	Fuel Type	Status	Capacity
F-1	None	Inactive	12.5 million gallons
F-2	JP-8	Active	12.5 million gallons
F-3	JP-8	Active	12.5 million gallons
F-4	JP-8	Active	12.5 million gallons
F-5	JP-8	Active	12.5 million gallons
F-6	JP-8	Active	12.5 million gallons
F-7	JP-5	Active	12.5 million gallons
F-8	JP-5	Active	12.5 million gallons
F-9	JP-5	Active	12.5 million gallons
F-10	JP-5	Active	12.5 million gallons
F-11	JP-5	Active	12.5 million gallons
F-12	JP-5	Active	12.5 million gallons
F-13	F-76	Active	12.5 million gallons
F-14	F-76	Active	12.5 million gallons
F-15	F-76	Active	12.5 million gallons
F-16	F-76	Active	12.5 million gallons
F-17	JP-5	Active	12.5 million gallons
F-18	JP-5	Active	12.5 million gallons
F-19	None	Inactive	12.5 million gallons
F-20	JP-5	Active	12.5 million gallons

F-76 Marine Diesel Fuel

JP-5 Jet Fuel Propellant-5

JP-8 Jet Fuel Propellant-8

#### 1.2 PHYSICAL SETTING

Climatological conditions in the area of the RHSF consist of warm to moderate temperatures and low to moderate rainfall. The RHSF is leeward of the prevailing northeasterly trade winds. The average annual precipitation is approximately 40 inches, which occurs mainly between November and April (State of Hawaii Department of Land and Natural Resources [DLNR], 1986). Annual pan evaporation is approximately 75 inches (DLNR, 1985). Average temperatures range from the low 60's to high 80's (degrees Fahrenheit) (Atlas of Hawaii, 1983).

Oahu consists of the eroded remnants of two shield volcanoes, Waianae and Koolau. The RHSF is located on the southwest flank of the Koolau volcanic shield. Lavas erupted during the shield-building phase of the volcano belong to the *Koolau Volcanic Series* (Stearns and Vaksvik, 1935). Following formation of the Koolau shield, a long period of volcanic quiescence occurred, during which the shield was deeply eroded. Following this erosional period, eruptive activity resumed. Lavas and pyroclastic material erupted during this period belong to the *Honolulu Volcanic Series* (Stearns and Vaksvik, 1935).

In the immediate area of the RHSF, Koolau Volcanic Series lavas dominate, although there are consolidated and unconsolidated non-calcareous deposits in the vicinity that consist of alluvium generated during erosion of the Koolau volcanic shield. South-southwest of the RHSF, and in isolated exposures to the west, are pyroclastic deposits formed during eruptions from three Honolulu Volcanic Series vents, Salt Lake, Aliamanu, and Makalapa (Stearns and Vaksvik, 1935). Based on established geology and records of the drilled wells (Stearns and Vaksvik, 1938), the RHSF is underlain by Koolau Volcanic Series basalts. The area of the RHSF is classified as *Rock Land*, where 25-90% of the land surface is covered by exposed rock and there are only shallow soils (Foote, et al., 1972).

Groundwater in Hawaii exists in two principal types of aquifers. The first and most important type, in terms of drinking water resources, is the basal aquifer. The basal aquifer exists as a lens of fresh water floating on and displacing seawater within the pore spaces, fractures, and voids of the basalt that forms the underlying mass of each Hawaiian island. In parts of Oahu, groundwater in the basal aquifer is confined by the overlying caprock and is under pressure. Waters that flow freely to the surface from wells that tap the basal aquifer are referred to as artesian.

The second type of aquifer is the caprock aquifer, which consists of various kinds of unconfined and semi-confined groundwater. Commonly, the caprock consists of a thick sequence of nearly impermeable clays, coral, and basalt, which separates the caprock aquifer from the basal aquifer. The impermeable nature of these materials and the artesian nature of the basal aquifer severely restrict the downward migration of groundwater from the upper caprock aquifer. In the area of the RHSF, there is no discernible caprock.

Groundwater in the area of the RHSF is part of the *Waimalu Aquifer System* of the *Pearl Harbor Aquifer Sector*. The aquifer is classified as a basal, unconfined, flank-type; and is currently used as a drinking water source. The aquifer is considered fresh with less than 250 milligrams per liter of chloride and is considered an irreplaceable resource with a high vulnerability to contamination (Mink and Lau, 1990).

The nearest drinking water supply well is DON Well 2254-01, located in the infiltration gallery within the RHSF lower tunnel. The DON Well 2254-01 is located approximately 2,400 feet hydraulically and topographically downgradient of the USTs (Figure 2).

#### 1.3 BACKGROUND

The RHSF was constructed by the U.S. Government in the early 1940s. Twenty USTs and a series of tunnels were constructed. The USTs were constructed of steel, and in the past have stored DON special fuel oil, DON distillate, aviation gasoline, and motor gasoline (Environet, 2010). The tanks currently contain JP-5, JP-8, and F-76. The fueling system is a self-contained underground unit that was installed into native rock comprised primarily of basalt with some interbedded tuffs and breccias (Environet, 2010). Each UST measures approximately 250 feet

in height and 100 feet in diameter. The upper domes of the tanks lie at a depth varying between 100 feet and 200 feet below ground surface [bgs].

In 1998, Earth Tech conducted a Phase II Remedial Investigation/Feasibility Study for the Oily Waste Disposal Facility located within the RHSF. The study included the installation of well OWDFMW01 (which was originally MW08) (Earth Tech, 1999).

In February 2001, the DON installed groundwater monitoring well RHMW01 to monitor for contamination in the basal aquifer beneath the RHSF. Well RHMW01 was installed approximately 100 feet below grade within the lower access tunnel. The depth to water was measured at 86 feet below the tunnel floor at the time of the well completion. In February 2001, a groundwater sample was collected from the well. Total Petroleum Hydrocarbons [TPH] and total lead were detected in the sample. Total lead was detected at a concentration above the DOH Tier 1 groundwater action level of 5.6  $\mu$ g/L (The Environmental Company, Inc. [TEC], 2009; DOH, 2000).

In 2005, the RHSF groundwater monitoring program was initiated. It involved routine groundwater sampling of well RHMW01 and sampling point RHMW2254-01. Samples were collected in February, June, September, and December of 2005. Lead was detected at concentrations above the DOH Tier 1 action level of 5.6 µg/L in samples collected in February and June. The samples collected in February and June were not filtered prior to analysis, whereas the samples collected in September and December were filtered prior to analysis. Since the samples collected in February and June were not filtered prior to analysis, the lead results were not considered appropriate for a risk assessment (TEC, 2008).

Between June and September 2005, TEC installed three additional groundwater monitoring wells (wells RHMW02, RHMW03, and RHMW04) (TEC, 2008). Well RHMW04 was installed hydraulically upgradient of the USTs to provide background geochemistry information for water moving through the basal aquifer beneath the RHSF. Wells RHMW02 and RHMW03 were installed approximately 125 feet below grade within the RHSF lower tunnel and well RHMW04 was installed to a depth of approximately 300 feet bgs outside of the RHSF tunnels. In September 2005, groundwater samples were collected from the three newly installed groundwater monitoring wells (wells RHMW02, RHMW03, and RHMW04) along with the existing well RHMW01 and sampling point RHMW2254-01. The contaminants of potential concern [COPCs] with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EAL.
- RHMW02 TPH-g, TPH-d, naphthalene, trichloroethylene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above the DOH EALs.
- RHMW03 TPH-d was detected at concentrations above the DOH EAL.

In 2006, TEC installed dedicated sampling pumps in the four wells (wells RHWM01, RHWM02, RHMW03, and RHWM04) and one sampling point (RHMW2254-01). In July and December of

2006, groundwater samples were collected from the four wells and the sampling point. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d and naphthalene were detected at concentrations above the DOH EALs.
- RHMW02 TPH-g, TPH-d, and naphthalene were detected at concentrations above the DOH EALs.
- RHMW03 TPH-d was detected at concentrations above the DOH EAL.

IN 2007, SSRBLs were established for TPH (4,500  $\mu$ g/L) and benzene (750  $\mu$ g/L) (TEC, 2007). Groundwater samples were collected from wells RHWM01, RHWM02, and RHMW03, and sampling point RHMW2254-01. Samples were collected in March, June, and September of 2007. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EAL, but below the SSRBL.
- RHMW02 TPH-g, TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above the DOH EALs. The TPH-d concentrations were below the SSRBL.
- RHMW03 TPH-d was detected at concentrations above the DOH EAL, but below the SSRBL.

In 2008, groundwater samples were collected from wells RHWM01, RHWM02, and RHMW03, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October of 2008. The COPCs with concentrations exceeding current DOH EALs are summarized below. In addition, a groundwater protection plan (TEC, 2008) was prepared.

- RHMW01 TPH-d was detected at concentrations above the DOH EAL, but below the SSRBL.
- RHMW02 TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above the DOH EALs. The TPH-d concentrations detected in October 2008 were also above the SSRBL.
- RHMW03 TPH-d was detected at concentrations above the DOH EAL, but below the SSRBL.

In April 2009, groundwater monitoring well RHMW05 was installed downgradient of the USTs, within the lower access tunnel between RHMW01 and RHMW2254-01. It was installed to identify the extent of contamination hydraulically downgradient of the USTs. Well RHMW05 was added to the quarterly groundwater sampling program. In 2009, quarterly groundwater samples were collect from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in February, May, July, and October of 2009. The COPCs with concentrations exceeding current DOH EALs are summarized below. In addition, the Groundwater Protection Plan was revised.

- RHMW01 TPH-d and 1-methylnaphthalene were detected at concentrations above the DOH EALs. The TPH-d concentrations were below the SSRBL.
- RHMW02 TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were
  detected at concentrations above the DOH EALs. The TPH-d concentrations were below the
  SSRBL.
- RHMW03 TPH-d was detected at a concentration above the DOH EAL, but below the SSRBL.
- RHMW05 TPH-d was detected at a concentration above the DOH EAL, but below the SSRBL.

In 2010, groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EAL, but below the SSRBI
- RHMW02 TPH-g, TPH-d, naphthalene, and 1-methylnaphthalene were detected at concentrations above the DOH EALs. The TPH-d concentrations were below the SSRBL.
- RHMW03 TPH-d was detected at a concentration above the DOH EAL, but below the SSRBL.
- RHMW05 TPH-d was detected at a concentration above the DOH EAL, but below the SSRBL.

In 2011, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. In a Fall 2011 update, the DOH EALs were revised. The drinking water toxicity EAL for TPH-d decreased from 210 to 190  $\mu$ g/L (DOH, 2011). The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EAL, but below the SSRBL.
- RHMW02 TPH-d, naphthalene, indeno[1,2,3-cd]pyrene, and 1-methylnaphthalene were
  detected at concentrations above the DOH EALs. The TPH-d concentrations were below the
  SSRBL.

In 2012, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in February, April, July, and November. The COPCs with concentrations exceeding current DOH EALs are summarized below.

 RHMW01 – TPH-d was detected at concentrations above the DOH EAL, but below the SSRBL.  RHMW02 – TPH-d, TPH-g, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above the DOH EALs. The TPH-d concentrations were below the SSRBL.

In 2013, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EAL, but below the SSRBL.
- RHMW02 TPH-d, TPH-g, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above the DOH EALs. The TPH-d concentrations were below the SSRBL.

In January 2014, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, RHMW05, and sampling point RHMW2254-01. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at a concentration above the DOH EAL, but below the SSRBL.
- RHMW02 TPH-d, naphthalene, and 1-methylnaphthalene were detected at concentrations above the DOH EALs. The TPH-d concentrations exceeded the SSRBL.

Between January and April 2014, additional groundwater sampling was conducted at wells RHMW01, RHMW02, RHMW05, and sampling point RHMW2254-01 in response to a suspected release from Tank 5. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EAL, but below the SSRBL.
- RHMW02 TPH-d, naphthalene, and 1-methylnaphthalene were detected at concentrations above the DOH EALs. The TPH-d concentrations were below the SSRBL.

In April 2014, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, RHMW05, and sampling point RHMW2254-01. The COPCs with concentrations exceeding current DOH EALs are summarized below.

 RHMW02 – TPH-d and naphthalene were detected at concentrations above both DOH EALs. 1-Methylnaphthalene was detected at concentrations above the DOH EAL for drinking water toxicity. The TPH-d concentrations were below the SSRBL.

In May and June 2014, additional groundwater sampling was conducted at wells RHMW01, RHMW02, RHMW05, and sampling point RHMW2254-01 in response to a reported release

from Tank 5. The COPCs with concentrations exceeding current DOH EALs are summarized below.

• RHMW02 – TPH-d, naphthalene, and 1-methylnaphthalene were detected at concentrations above the DOH EALs. The TPH-d concentrations were below the SSRBL.

#### 1.3.1 Previous Reports

The following groundwater monitoring reports were previously submitted to the DOH:

- Groundwater Sampling Report, First Quarter 2005 (submitted April 2005).
- Groundwater Sampling Report, Second Quarter 2005 (submitted August 2005).
- 3. Groundwater Sampling Report, Third Quarter 2005 (submitted November 2005).
- Groundwater Sampling Report, Fourth Quarter 2005 (submitted February 2006).
- 5. Groundwater Monitoring Results, July 2006 (submitted September 2006).
- 6. Groundwater Monitoring Results, December 2006 (submitted January 2007).
- 7. Groundwater Monitoring Results, March 2007 (submitted May 2007).
- 8. Groundwater Monitoring Results, June 2007 (submitted August 2007).
- 9. Groundwater Monitoring Results, September 2007 (submitted October 2007).
- 10. Groundwater Monitoring Report, January 2008 (submitted March 2008).
- 11. Groundwater Monitoring Report, April 2008 (submitted May 2008).
- 12. Groundwater Monitoring Report, July 2008 (submitted October 2008).
- 13. Groundwater Monitoring Report, October and December 2008 (submitted February 2009).
- 14. Groundwater Monitoring Report, February 2009 (submitted May 2009).
- 15. Groundwater Monitoring Report, May 2009 (submitted July 2009).
- 16. Groundwater Monitoring Report, July 2009 (submitted September 2009).
- 17. Groundwater Monitoring Report, October 2009 (submitted December 2009).
- 18. Groundwater Monitoring Report, January, February, and March 2010 (submitted April 2010).
- 19. Groundwater Monitoring Report, April 2010 (submitted May 2010).

- 20. Groundwater Monitoring Report, July 2010 (submitted August 2010).
- 21. Groundwater Monitoring Report, October 2010 (submitted December 2010).
- 22. Groundwater Monitoring Report, January 2011 (submitted March 2011).
- 23. Groundwater Monitoring Report, April 2011 (submitted June 2011).
- 24. Groundwater Monitoring Report, July 2011 (submitted September 2011).
- 25. Groundwater Monitoring Report, October 2011 (submitted December 2011).
- 26. Groundwater Monitoring Report, January-February 2012 (submitted March 2012).
- 27. Groundwater Monitoring Report, April 2012 (Submitted July 2012).
- 28. Groundwater Monitoring Report, October 2012 (Submitted January 2013).
- 29. Groundwater Monitoring Report, January 2013 (Submitted April 2013).
- 30. Groundwater Monitoring Report, April 2013 (Submitted July 2013).
- 31. Groundwater Monitoring Report, July 2013 (Submitted September 2013).
- 32. Groundwater Monitoring Report, October 2013 (Submitted January 2014).
- 33. Groundwater Sampling Report for Additional Sampling, January 2014 (submitted January 2014).
- 34. Groundwater Monitoring Report, January 2014 (Submitted April 2014).
- 35. Groundwater Sampling Report for Tank 5 Release Response on March 5 and 6, 2014 (submitted March 2014).
- 36. Groundwater Sampling Report for Tank 5 Release Response on March 10, 2014 (submitted March 2014).
- 37. Groundwater Sampling Report for Tank 5 Release Response on March 25 and 26, 2014 (submitted April 2014).
- 38. Groundwater Sampling Report for Tank 5 Release Response on April 7, 2014 (submitted April 2014).
- 39. Groundwater Monitoring Report, April 2014 (Submitted June 2014).
- 40. Groundwater Sampling Report for Tank 5 Release Response on May 27 and 28, 2014 (submitted June 2014).

41. Groundwater Sampling Report for Tank 5 Release Response on June 23 and 24, 2014 (submitted July 2014).

#### **SECTION 2 – GROUNDWATER SAMPLING**

On July 21 and 22, 2014, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01). The samples were collected in accordance with the 2012 WP/SAP. The WP/SAP is consistent with DOH UST release response requirements (DOH, 2000); DoN Procedure I-C-3, *Monitoring Well Sampling* (DoN, 2007); and the RHSF Groundwater Protection Plan (TEC, 2008). Prior to purging and sampling, the depth to groundwater and the depth to the bottoms of the wells were measured using a Geotech oil/water interface probe. No measurable product, sheen, or petroleum hydrocarbon odor was detected in any of the wells with the exception of RHMW02. A moderate petroleum hydrocarbon odor was noticed when collecting groundwater samples from RHMW02.

#### 2.1 GROUNDWATER SAMPLING

Prior to collecting groundwater samples, the monitoring wells were purged of standing water in the well casings. Each well contains a dedicated bladder pump which was used to purge the well and to collect samples. To operate the pump, a portable air compressor with an in-line filter was connected to a QED MP50 MicroPurge® Basics Controller box, which was then connected to the pump. The compressor was turned on to power the pump and the controller was used to adjust the pumping rate to less than one liter of water per minute.

Water quality parameters were monitored on a periodic basis during well purging. The water quality parameters that were measured included hydrogen activity [pH], temperature, conductivity, dissolved oxygen, and oxidation reduction potential. The water quality parameters were evaluated to assess whether the natural characteristics of the aquifer formation water were present within the monitoring wells before collecting the samples. At least four readings were collected during the purging process. Purging was considered complete when at least three consecutive water quality measurements stabilized within approximately 10%. The readings were recorded on groundwater monitoring logs. The groundwater monitoring logs are included in Appendix A. In addition, field notes were taken to document the sampling event. The field notes are included in Appendix B.

When the water quality parameters stabilized, groundwater samples were collected from the wells using the bladder pumps. The groundwater samples were collected no more than two hours after purging was completed to decrease groundwater interaction with the monitoring well casing and atmosphere. Prior to collecting the sample, the water level in the monitoring wells was measured and recorded to ensure that excessive drawn down had not occurred. The groundwater samples were collected at flow rates of approximately 0.17 to 0.5 liters per minute. Samples collected for dissolved lead analysis were filtered in the field using new 0.45 micron filters.

#### 2.2 ANALYTICAL RESULTS

The samples were analyzed for TPH-d using U.S. Environmental Protection Agency [EPA] Method 8015M, TPH-g and Volatile Organic Compounds [VOCs] using EPA Method 8260B, PAHs using EPA Method 8270C SIM, dissolved lead using EPA Method 6020, and total lead using EPA Method 200.8. The sample collected from sampling point RHMW2254-01 was analyzed for total lead (unfiltered) as DON Well 2254-01 is a drinking water supply well. The analytical results are described below and summarized in Table 2.1. A copy of the laboratory report is included in Appendix C. Analytical results were compared to the DOH EALs for drinking water toxicity and gross contamination and to the SSRBLs established in the RHSF Final Groundwater Protection Plan (TEC, 2008).

- RHMW01 TPH-d (67 μg/L) was the only analyte detected. The TPH-d concentration detected did not exceed the DOH EALs or the SSRBL.
- RHMW02 TPH-d (1,200 and 1,300 μg/L), TPH-g (48 and 49 μg/L), xylenes (0.36 and 0.33 μg/L), acenaphthene (0.52 and 0.50 μg/L), fluorene (0.24 and 0.23 μg/L), 1-methylnaphthalene (25 and 26 μg/L), 2-methylnaphthalene (20 and 22 μg/L), and naphthalene (71 and 76 μg/L) were detected. TPH-d, 1-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EALs for both drinking water toxicity and gross contamination. However, the TPH-d concentrations did not exceed the SSRBL. 2-Methylnaphthalene was detected at concentrations above the DOH EAL for gross contamination.
- RHMW03 TPH-d (37 μg/L) was the only analyte detected. None of the chemical constituents analyzed for were detected at concentrations above he DOH EALs or the SSRBL.
- RHMW05 None of the chemical constituents analyzed for were detected.
- RHMW2254-01 None of the chemical constituents analyzed for were detected.

Acetone (6.5  $\mu$ g/L) was detected in the trip blank submitted on July 21, 2014. Acetone was not detected in any of the groundwater samples, and this is not likely to affect data usability.

#### 2.3 GROUNDWATER CONTAMINANT TRENDS

The historical groundwater contaminant concentration trends for COPCs that exceeded the DOH EALs or SSRBLs are illustrated in Appendix D. A summary of groundwater contaminant trends is provided below.

 RHMW01 – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW01. TPH-d has historically been detected at concentrations above the DOH EAL for both drinking water toxicity and gross contamination. TPH-d concentrations have shown an overall decreasing trend from a high of 1,500  $\mu$ g/L in February 2005. The TPH-d concentration was last above the DOH EAL for gross contamination in a sample collected on April 7, 2014.

- RHMW02 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW02. TPH-g, TPH-d, trichloroethylene, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene have historically been detected at concentrations above the DOH EALs. TPH-d concentrations increased sharply from 2,400 µg/L to 5,100 µg/L during the additional sampling event on January 15, 2014, exceeding the SSRBL of 4,500 µg/L. During this July 2014 event, TPH-d concentrations decreased to an average of both 1,250 µg/L but remained above DOH EALs. The concentrations 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene increased to their highest levels since January 2013. The concentrations of TPH-g remained below the DOH EALs for gross contamination and drinking water toxicity. Trichloroethylene has not been detected in RHMW02 since September 2005.
- RHMW03 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW03. TPH-d has historically been detected at concentrations above the DOH EALs; however, it has not been detected at concentrations above the DOH EALs since October 2010.
- RHMW05 No COPCs were detected during this round of quarterly sampling. TPH-d has
  historically been detected in RHMW05 at concentrations above the DOH EAL for both
  drinking water toxicity and gross contamination; however, it has not been detected at
  concentrations above the DOH EALs since January 2010.
- RHMW2254-01 No COPCs were detected during this round of quarterly sampling. TPH-d was last detected in RHMW2254-01 at a concentration above the DOH EAL for gross contamination in January 2008. Although the method reporting limits for TPH-d exceeded one or both DOH EALs between May 2009 and July 2010, TPH-d has not been detected at concentrations above the DOH EALs since January 2008.

Historical groundwater contaminant trends for concentrations above the DOH EALs for both drinking water toxicity and gross contamination are illustrated in Appendix D.

#### 2.4 WASTE DISPOSAL

The purged groundwater and decontamination water generated during sampling of the inside tunnel wells were stored in a 55-gallon drum along with the purged water and decontamination water from the outside tunnel wells. The drum is currently stored onsite at ADIT 3 on top of a secondary containment spill pallet and covered by a tarp. There is a non-hazardous label affixed to the drum with all pertinent information relating to its generation. The drum will be used for future sampling events and will be properly disposed of once it has been filled.

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#### **TABLE 2.1** Analytical Results for Groundwater Sampling (July 21 and 22, 2014) Red Hill Bulk Fuel Storage Facility **July 2014 Quarterly Monitoring Report**

		DOH EALs				2254-01 (ES	S107)				M01 (ES10		iy 14.01.1	RHMW02 (ES104)						DHWM	//03 (ES1	06)		RHMW05 (ES108)				
Method	Chemical	Drinking Water	Gross	Results	Q	LOQ	LOD	DL	Results	Q	LOQ	LOD	DL	Results	Q	LOQ	LOD	DL	Results	Q	LOQ	LOD	DL	Results	Q	LOQ	LOD	DL
EPA 8015B	TPH-d	Toxicity 190	Contamination 100	N.D.	U	25	12	7.4	67	HD	25	10	7.4	1,200	HD	25	10	7.4	37	HD	25	10	7.4	N.D.	U	25	12	7.4
EPA 8260B	TPH-q	100	100	N.D.	U	50	30	26	N.D.	U	50	30	26	48	J	50	30	26	N.D.	U	50	30	26	N.D.	U	50	30	26
	Acenaphthene	370	20	N.D.	Ū	0.2	0.048	0.020	N.D.	U	0.2	0.051	0.021	0.52	Ť	0.2	0.048	0.021	N.D.	Ü	0.2	0.047	0.020	N.D.	U	0.2	0.049	0.021
	Acenaphthylene	240	2,000	N.D.	U	0.2	0.048	0.018	N.D.	U	0.2	0.051	0.018	N.D.	U	0.2	0.048	0.018	N.D.	U	0.2	0.047	0.018	N.D.	U	0.2	0.049	0.018
	Anthracene	1,800	22	N.D.	U	0.2	0.048	0.034	N.D.	U	0.2	0.051	0.034	N.D.	U	0.2	0.048	0.035	N.D.	U	0.2	0.047	0.034	N.D.	U	0.2	0.049	0.035
	Benzo[a]anthracene	0.092	4.7	N.D.	U	0.2	0.048	0.023	N.D.	U	0.2	0.051	0.024	N.D.	U	0.2	0.048	0.024	N.D.	U	0.2	0.047	0.023	N.D.	U	0.2	0.049	0.024
	Benzo[g,h,i]perylene	1,500	0.13	N.D.	U	0.2	0.048	0.022	N.D.	U	0.2	0.051	0.022	N.D.	U	0.2	0.048	0.022	N.D.	U	0.2	0.047	0.022	N.D.	U	0.2	0.049	0.022
	Benzo[a]pyrene	0.2	0.81	N.D.	U	0.2	0.048	0.036	N.D.	U	0.2	0.051	0.037	N.D.	U	0.2	0.048	0.037	N.D.	U	0.2	0.047	0.036	N.D.	U	0.2	0.049	0.037
	Benzo[b]fluoranthene	0.092	0.75	N.D.	U	0.2	0.048	0.025	N.D.	U	0.2	0.051	0.025	N.D.	U	0.2	0.048	0.025	N.D.	U	0.2	0.047	0.025	N.D.	U	0.2	0.049	0.025
	Benzo[k]fluoranthene	0.92	0.4	N.D.	U	0.2	0.048	0.023	N.D.	U	0.2	0.051	0.024	N.D.	U	0.2	0.048	0.024	N.D.	U	0.2	0.047	0.023	N.D.	U	0.2	0.049	0.024
EPA 8270C	Chrysene	9.2 0.0092	0.52	N.D.	U	0.2	0.048	0.019	N.D. N.D.	U	0.2	0.051 0.051	0.019 0.027	N.D. N.D.	U	0.2	0.048	0.019 0.027	N.D. N.D.	U	0.2	0.047 0.047	0.019	N.D. N.D.	U	0.2	0.049 0.049	0.019 0.027
	Dibenzo[a,h]anthracene   Fluoranthene	1,500	130	N.D.	U	0.2	0.048	0.026	N.D.	U	0.2	0.051	0.027	N.D.	U	0.2	0.048	0.027	N.D.	U	0.2	0.047	0.026	N.D.	U	0.2	0.049	0.027
	Fluorene	240	950	N.D.	U	0.2	0.048	0.024	N.D.	U	0.2	0.051	0.027	0.24	-	0.2	0.048	0.025	N.D.	U	0.2	0.047	0.024	N.D.	U	0.2	0.049	0.025
	Indeno[1,2,3-cd]pyrene	0.092	0.095	N.D.	U	0.2	0.048	0.024	N.D.	Ü	0.2	0.051	0.023	N.D.	U	0.2	0.048	0.023	N.D.	U	0.2	0.047	0.024	N.D.	U	0.2	0.049	0.023
ĺ	1-Methylnaphthalene	4.7	10	N.D.	U	0.2	0.048	0.028	N.D.	U	0.2	0.051	0.029	25	<u> </u>	0.2	0.96	0.029	N.D.	U	0.2	0.047	0.028	N.D.	U	0.2	0.049	0.029
ĺ	2-Methylnaphthalene	24	10	N.D.	Ü	0.2	0.048	0.026	N.D.	Ü	0.2	0.051	0.027	20		0.2	0.96	0.027	N.D.	Ü	0.2	0.047	0.027	N.D.	Ü	0.2	0.049	0.027
	Naphthalene	17	21	N.D.	U	0.2	0.048	0.023	N.D.	U	0.2	0.051	0.023	71		2.0	0.96	0.23	N.D.	U	0.2	0.047	0.023	N.D.	U	0.2	0.049	0.023
ĺ	Phenanthrene	240	410	N.D.	U	0.2	0.048	0.030	N.D.	U	0.2	0.051	0.031	N.D.	U	0.2	0.048	0.031	N.D.	U	0.2	0.047	0.030	N.D.	U	0.2	0.049	0.031
	Pyrene	180	68	N.D.	U	0.2	0.048	0.024	N.D.	U	0.2	0.051	0.025	N.D.	U	0.2	0.048	0.025	N.D.	U	0.2	0.047	0.024	N.D.	U	0.2	0.049	0.025
I	1,1,1,2-Tetrachloroethane	0.52	50,000	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4
ĺ	1,1,2,2-Tetrachloroethane <sup>1</sup>	0.067	500	N.D.	U	1	0.5	0.41	N.D.	U	1 -	0.5	0.41	N.D.	U	1	0.5	0.41	N.D.	U	1 -	0.5	0.41	N.D.	U	1 -	0.5	0.41
ĺ	1,1,1-Trichloroethane	200	970	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3	N.D.	U	5	0.5	0.3
	1,1,2-Trichloroethane	5	50,000	N.D.	U	1	0.5	0.38	N.D.	U	1 5	0.5	0.38	N.D.	U	1	0.5	0.38	N.D.	U	1	0.5	0.38	N.D.	U	1	0.5	0.38
	1,1-Dichloroethane	2.4	50,000	N.D.	U	5	0.5	0.28	N.D.	U	5	0.5	0.28	N.D.	U	5 1	0.5	0.28	N.D.	U	5	0.5	0.28	N.D.	U	5 1	0.5	0.28
	1,1-Dichloroethylene 1,2,3-Trichloropropane <sup>1</sup>	0.6	1,500 50,000	N.D.	U	5	0.5 1	0.43 0.64	N.D. N.D.	U	5	0.5 1	0.43	N.D. N.D.	U	5	0.5	0.43	N.D. N.D.	U	5	0.5	0.43	N.D. N.D.	U	5	0.5	0.43 0.64
	1,2,4-Trichlorobenzene	70	3,000	N.D.	U	5	1	0.5	N.D.	U	5	1	0.04	N.D.	U	5	1	0.04	N.D.	U	5	1	0.5	N.D.	U	5	1	0.5
	1,2-Dibromo-3- chloropropane <sup>1</sup>	0.04	10	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2	N.D.	U	10	2	1.2
	1,2-Dibromoethane <sup>1</sup>	0.04	50,000	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24
	1,2-Dichlorobenzene	600	10	N.D.	Ü	1	0.5	0.46	N.D.	Ü	1	0.5	0.46	N.D.	Ü	1	0.5	0.46	N.D.	Ü	1	0.5	0.46	N.D.	U	1	0.5	0.46
	1,2-Dichloroethane <sup>1</sup>	0.15	7,000	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24
	1,2-Dichloropropane	5	10	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42
	1,3-Dichlorobenzene	180	5	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4
	1,3-Dichloropropene (total of cis/trans) 1	0.43	50,000	N.D.	J	1	0.5	0.25	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25
	1,4-Dichlorobenzene	75	5	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43
	Acetone	22,000	20,000	N.D.	U,ICH	20	10	6	N.D.	U,ICH	20	10	6	N.D.	U,ICH	20	10	6	N.D.	U,ICH	20	10	6	N.D.	U,ICH	20	10	6
	Benzene	5	170	N.D.	U	1	0.5	0.14	N.D.	U	1	0.5	0.14	N.D.	U	1	0.5	0.14	N.D.	U	1	0.5	0.14	N.D.	U	1	0.5	0.14
	Bromodichloromethane <sup>1</sup>	0.12	50,000	N.D.	U	5	0.5	0.21	N.D.	U	5	0.5	0.21	N.D.	U	5	0.5	0.21	N.D.	U	5	0.5	0.21	N.D.	U	5	0.5	0.21
	Bromoform	80	510	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5
EPA 8260B	Bromomethane	8.7	50,000	N.D.	U	20	5	3.9	N.D.	U	20	5	3.9	N.D.	U	20	5	3.9	N.D.	U	20	5	3.9	N.D.	U	20	5	3.9
	Carbon Tetrachloride	5	520	N.D.	U	1	0.5	0.23	N.D.	U	1 5	0.5	0.23	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23
ĺ	Chlorobenzene Chloroethane	100 21,000	50 16	N.D. N.D.	U	5 10	0.5 5	0.17 2.3	N.D. N.D.	U	5 10	0.5 5	0.17 2.3	N.D. N.D.	U	5 10	0.5 5	0.17 2.3	N.D. N.D.	U	5 10	0.5 5	0.17 2.3	N.D. N.D.	U	5 10	0.5 5	0.17 2.3
ĺ	Chloroform	70	2,400	N.D.	U	5	0.5	0.46	N.D.	U	5	0.5	0.46	N.D.	U	5	0.5	0.46	N.D.	U	5	0.5	0.46	N.D.	U	5	0.5	0.46
	Chloromethane <sup>1</sup>	1.8	50,000	N.D.	U, IJ	10	2	1.8	N.D.	U, IJ	10	2	1.8	N.D.	U, IJ	10	2	1.8	N.D.	U, IJ	10	2	1.8	N.D.	U, IJ	10	2	1.8
	cis-1,2-Dichloroethylene	70	50,000	N.D.	U	1	0.5	0.48	N.D.	U	1	0.5	0.48	N.D.	U	1	0.5	0.48	N.D.	U	1	0.5	0.48	N.D.	U	1	0.5	0.48
l	Dibromochloromethane <sup>1</sup>	0.16	50,000	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25
	Ethylbenzene	700	30	N.D.	U	1	0.5	0.14	N.D.	Ü	1	0.5	0.14	N.D.	J	1	0.5	0.14	N.D.	J	1	0.5	0.14	N.D.	U	1	0.5	0.14
	Hexachlorobutadiene	0.86	6	N.D.	U	1	0.5	0.32	N.D.	U	1	0.5	0.32	N.D.	U	1	0.5	0.32	N.D.	U	1	0.5	0.32	N.D.	U	1	0.5	0.32
	Methyl ethyl ketone (2- Butanone)	7,100	8,400	N.D.	J	10	5.0	2.2	N.D.	U	10	5.0	2.2	N.D.	U	10	5.0	2.2	N.D.	U	10	5.0	2.2	N.D.	U	10	5.0	2.2
	Methyl isobutyl ketone (4- Methyl-2-Pentanone)	2,000	1300	N.D.	U	10	5.0	4.4	N.D.	U	10	5.0	4.4	N.D.	U	10	5.0	4.4	N.D.	U	10	5.0	4.4	N.D.	U	10	5.0	4.4
	Methyl tert-butyl Ether	12	5	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31
	Methylene chloride	4.8	9,100	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64
	Styrene	100	10	N.D.	U	1	0.5	0.17	N.D.	U	1	0.5	0.17	N.D.	U	1	0.5	0.17	N.D.	U	1	0.5	0.17	N.D.	U	1	0.5	0.17
	Tetrachloroethylene	5	170	N.D.	U	5	0.5	0.39	N.D.	U	5	0.5	0.39	N.D.	U	5	0.5	0.39	N.D.	U	5	0.5	0.39	N.D.	U	5	0.5	0.39
	Toluene	1,000	40	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24
	trans-1,2- Dichloroethylene	100	260	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	N.D.	U	1 1	0.5	0.37	N.D.	U	1	0.5	0.37
	Trichloroethylene	5	310	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37
	Vinyl chloride  Xylenes	10,000	3,400 20	N.D. N.D.	U	1 11	0.5 1.5	0.3	N.D. N.D.	U	1	0.5	0.3	N.D. 0.36	U	1	0.5 0.5	0.3	N.D.	U U	1	0.5	0.3	N.D.	U	1 11	0.5 1.5	0.3 0.23
EPA 6020	Dissolved Lead	10,000	50,000	N.D.	-	- 11	1.5	0.23	N.D.	U	1	0.5	0.23	0.36 N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23	N.D.	U	11	0.2	0.23
EPA 200.8	Total Lead	15	50,000	N.D.	U	1.0	-	0.0898	14.0.	-	-	-	0.0090	IN.D.	-	-	-	0.0090	N.D.	-	-	- 0.2	0.0096	N.D.	-	-	-	-
21 // 200.0	. Star Edua	10	55,000	11.0.	U	1.0	_	0.0000	1	-	1 -	-		_		_	_		1	•			1 -		•		•	

The data are in micrograms per liter ( $\mu$ g/L). Shaded values exceeded the DOH EALs.

1 LOD for this analyte exceeds the DOH EAL

- Not Analyzed

DL DOH EALS EPA HD ICH ICJ

Not Analyzed
This sample was analyzed by EPA Method 200.8 and therefore does not have an LOD
Detection Limit or Method Detection Limit (MDL)
DOH Tier 1 Environmental Action Levels for groundwater where groundwater is a current drinking water source and surface water is greater than 150 meters from the site (DOH, Fall 2011).
Environmental Protection Agency
The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
Initial calibration verification recovery is above the control limit for this analyte.
Initial calibration verification recovery is below the control limit for this analyte.

Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated. Limit of Detection Limit of Quantitation
Not Detected Qualifiers
Total Petroleum Hydrocarbons as diesel

LOD LOQ N.D. Q TPH-d TPH-g U

Total Petroleum Hydrocarbons as gasoline
Undetected at DL and is reported as less than the LOD.

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#### **SECTION 3 – DATA QUALITY ASSESSMENT**

A data quality assessment, which consists of a review of the overall groundwater sample collection and analysis process, was performed in order to determine whether the analytical data generated met the quality objectives for the project. The data quality assessment was performed in accordance with the approved WP/SAP (ESI, 2012). The field quality control program consisted of standardized sample collection and management procedures, and the collection of field duplicate samples, matrix spike samples, and trip blank samples. The laboratory quality assurance program consisted of the use of standard analytical methods and the preparation and analyses of Matrix Spike [MS]/Matrix Spike Duplicate [MSD] samples, surrogate spikes, blanks, and Laboratory Control Samples [LCSs]/Laboratory Control Sample Duplicates [LCSDs].

#### 3.1 DATA VALIDATION AND ASSESSMENT

The objective of data validation is to provide data of known quality for project decisions. Data quality is judged in terms of Precision, Accuracy, Representativeness, Completeness, Comparability, and Sensitivity [PARCCS]. A number of factors may affect the quality of data, including: sample collection methods, sample analysis methods, and adherence to established procedures for sample collection, preservation, management, shipment, and analysis.

#### Precision

Precision is defined as the reproducibility of replicate measurements. Precision is evaluated by Relative Percentage Difference [RPD] of field duplicates, LCS/LCSDs, and MS/MSD results. Field duplicate and MS/MSD samples were collected at a rate of approximately 10% of primary samples. Field duplicates were sent to the laboratory along with the primary samples.

The RPDs of detected analytes for the primary and field duplicate samples (ES104 and ES105) are provided in Table 3.1. A precision of less than 50% for duplicate pairs is required by the DoN Project Procedures Manual to be considered acceptable (DoN 2007). All duplicate RPDs are less than the acceptable maximum, except for lead. An RPD of 200% was assigned to the ES104 and ES105 lead results because lead was detected in only one of the replicate samples from the duplicate pair. The lead result for sample ES105 was below the LOQ; consequently, the RPD exceedance signified the anticipated decrease in precision below the LOQ, but was not indicative of a QC issue.

Lead was detected in only the duplicate sample ES105, and was assigned an RPD of 200%. However, the concentration detected was below the limit of quantitation [LOQ] and well below the DOH EALs. Therefore, it is unlikely to have an effect on data usability, and the data precision is considered acceptable. In addition, all RPDs for MS/MSD and LCS/LCSD pairs were also below the control limit.

#### Accuracy

Accuracy is defined as the degree of conformity of a measurement to a standard or true value. Accuracy is evaluated through measurement of the percent recovery of an analyte in a reference standard or spiked sample. Accuracy limits for surrogates, laboratory control spike, MS, and MSD samples are established by the individual laboratory.

Between July 2006 and July 2010, naphthalene was analyzed for by both EPA Methods 8260B and 8270C, and both results were reported. In September 2005 and in all data beginning in October 2010, only results using EPA Method 8270C were reported. Naphthalene has historically only been detected at concentrations above the DOH EALs in well RHMW02. In this well, concentrations of naphthalene detected in each sample by EPA Method 8260B were generally two to three times higher than those detected by EPA Method 8270C. We assume this is due to the better preservation of VOCs associated with the use of EPA Method 8260B. This suggests that the naphthalene results provided by EPA Method 8270C may be biased low. Since October 2012, naphthalene concentrations in RHMW02 have exceeded DOH EALs for both gross contamination and drinking water toxicity. The naphthalene concentration detected in July 2012 (17 µg/L) was equal to the DOH EAL for drinking water (17 µg/L) but below the DOH EAL for gross contamination (21 µg/L); it is possible that accounting for the low bias, the actual naphthalene concentration detected in July exceeded both EALs. Naphthalene concentrations between April 2011 and April 2012 were all an order of magnitude below both EALs, and it is unlikely that naphthalene concentrations in these samples exceeded the EALs, even after accounting for a potential low bias.

Results for TPH-d in samples ES103 through ES106 were flagged "HD." The laboratory indicated a mismatch between the calibration standard and the TPH-d chromatographic profile. Mismatches of this type are not uncommon. Even though chromatograms are not part of the standard laboratory package, ESI was able to review the chromatograms from RHMW02 dating back to October 2012. The chromatograms of groundwater samples from RHMW02 did not significantly differ between each event, but did not match a standard chromatogram of JP-8 in groundwater.

For this July 2014 sampling event, all of the LCS and surrogate spike recoveries for analyzed constituents were within acceptable percent recovery limits, except for the LCS percent recovery for acetone, which was above the recovery limit. This indicates acetone results may be biased high; however, acetone was only detected in the trip blank and not in any of the groundwater samples.

The MS and/or MSD recoveries were above the control limits for acetone and lead and the associated sample results may be biased high. Acetone was not detected in any of the groundwater samples and lead was only detected in ES105, at a concentration below the LOQ. Naphthalene, 2-methylnaphthalene and 1-methylnaphthalene concentrations for ES104, the primary sample on which the MS/MSD were performed, were significantly higher than the added spike concentration, which prevented an accurate evaluation of the MS/MSD recovery for these analytes.

All other MS/MSD recoveries were within acceptable recovery limits; therefore, the data accuracy for this monitoring event is considered acceptable.

#### Representativeness

Representativeness is the degree that data accurately and precisely represents a characteristic of a population, parameter variations at a sampling point, or an environmental condition. Representativeness was achieved by conducting sampling in accordance with the sample collection procedures described in the project WP/SAP, including standardized sample collection methods (ESI, 2012).

Representativeness is also evaluated through the compliance with the standardized sample holding time and sample preservation methods, and through the analysis of blank samples, including method blank and trip blank samples. For this sampling event, all sample holding time and sample preservation were consistent with EPA guidance.

For this sampling event, one trip blank was included in every cooler containing samples for VOC and TPH-g analysis to assess the potential for contamination during sample transport. Two trip blanks were collected. Acetone was the only COPC detected, and in only one of the trip blanks; however, acetone was not detected in any of the groundwater samples, excluding the possibility that project samples were contaminated with acetone during handling or transport. Based on the assessment of representativeness, the groundwater sample data are considered representative of the groundwater quality on site. The trip blank results are provided in Table 3.1.

#### Completeness

Completeness is defined as the overall percentage of valid analytical results (including estimated results) compared to the total number of analytical results reported by the analytical laboratory. No data were rejected for this project, and therefore the completeness goal for this project (90%), was successfully met.

#### Comparability

Comparability expresses the confidence with which one data set can be compared to another data set. Comparability can be related to accuracy and precision because these quantities are measures of data reliability. Data with acceptable precision and accuracy are considered comparable if collection techniques, analytical procedures, methods and reporting are equivalent.

As noted above, between July 2006 and July 2010, naphthalene was analyzed for using both EPA Methods 8260B and 8270C, and in September 2005 and between October 2010 and the most recent event, only results using EPA Method 8270C were reported. In general, EPA Method 8260B resulted in higher, and as discussed above, likely more accurate, results than EPA Method 8270C. However, for the sake of comparability with results from recent events, EPA Method 8270C was used for naphthalene analysis in this event. Consequently, the low bias associated with Method 8270C should be considered when making project decisions.

All project samples for TPH-g analysis through July 2010 were analyzed by EPA Method 8015; beginning in October 2010, EPA Method 8260B was used. There was no event where both methods were used; consequently, there is no way to directly compare the results obtained by method and to assess potential bias. However, there is no reason to believe that using either method should bias the data, and the TPH-g data for all events should be comparable.

Other than the naphthalene bias discussed above, no issues with comparability were identified. The results are considered comparable within this data set and with the data collected from recent sampling events.

#### Sensitivity

The LOQs are established by the laboratory based on the limits of detection [LODs] or instrument detection limits, historical data, and EPA limits established for the various methods. The LOQs for samples may require adjustment by the laboratory due to matrix interference or if high levels of target analytes necessitate dilution before analysis. Matrix interference and sample dilutions have the effect of decreasing sensitivity and increasing the LOQs. Laboratory LODs and LOQs for several analytes (EPA Methods 8260 and 8270) for this event differed from the LODs and LOQs in the WP/SAP because the laboratory updates them quarterly and in some cases, dilution was necessary due to the presence of high concentrations of analytes.

For this event, LODs and LOQs for several analytes were greater than the DOH EALs (as stated in the WP/SAP), and therefore it would not be possible to detect the analytes at concentrations greater than the DOH EALs but below the LODs or LOQs. The lack of the required sensitivity should be considered when making project decisions. The affected analytes for this monitoring event are 1,2,3-trichloropropane, 1,2-dibromo-3-chloropropane, 1,2-dibromoethane, 1,2-dichloroethane, 1,3-dichloropropene, bromodichloromethane, chloromethane, dibromochloromethane, 1,1,2,2-tetrachloroethane, and dibenzo[a,h]anthracene.

#### 3.2 DATA ASSESSMENT AND USABILITY CONCLUSIONS

The PARCCS criteria were evaluated, and with a few exceptions, all criteria were met. The data assessment concludes that all data generated during this event are usable for the intended purpose.

Contract No. N62742-12-D-1853 Contract Task Order 0002

TABLE 3.1 Quality Control Results for Groundwater Sampling (July 21 and 22, 2014)

Red Hill Bulk Fuel Storage Facility

July 2014 Quarterly Monitoring Report

		DOH E		RHMW02 (ES104) RHMW02 (ES1										ES Trip (7/21/2014)						ES Trip (7/22/2014)				
	Chemical Constituent	DONE	LALS		I KITIW	W02 (L3	104)			KIIIVIVV	02 (23103	) (501)		RPD		LOT	iip (1/21/2	1014)						
Method		Drinking Water Toxicity	Gross Contamination	Result	Q	LOQ	LOD	DL	Result	Q	LOQ	LOD	DL	Duplicate (%)	Result	Q	LOQ	LOD	DL	Result	Q	LOQ	LOD	DL
EPA 8015B	TPH-d	190	100	1,200	HD	25	10	7.4	1,300	HD	25	10	7.4	8.00	-	-	-	-	-	-	-	-	-	-
EPA 8260B	TPH-g	100	100	48	J	50	30	26	49	J	50	30	26	2.06	N.D.	U	50	30	26	N.D.	U	50	30	26
	Acenaphthylone	370 240	20 2.000	0.52 N.D.	U	0.2	0.048 0.048	0.021 0.018	0.50 N.D.	U	0.2	0.05	0.021 0.018	3.92 NA	-	-	-	-	-	-	-	-	-	-
	Acenaphthylene Anthracene	1,800	22	N.D.	U	0.2	0.048	0.018	N.D.	U	0.2	0.05	0.018	NA NA	-	-	-	-	-	-	-	-	-	-
	Benzo[a]anthracene	0.092	4.7	N.D.	Ü	0.2	0.048	0.024	N.D.	Ü	0.2	0.05	0.024	NA	-	-	-	-	-	-	-	-	-	-
	Benzo[g,h,i]perylene	1,500	0.13	N.D.	U	0.2	0.048	0.022	N.D.	U	0.2	0.05	0.022	NA	-	-	-	-	-	-	-	-	-	-
	Benzo[a]pyrene	0.2	0.81	N.D.	U	0.2	0.048	0.037	N.D.	U	0.2	0.05	0.036	NA	-	-	-	-	-	-	-	-	-	-
	Benzo[b]fluoranthene	0.092	0.75	N.D.	U	0.2	0.048	0.025	N.D.	U	0.2	0.05	0.025	NA NA	-	-	-	-	-	-	-	-	-	-
	Benzo[k]fluoranthene Chrysene	0.92 9.2	0.4	N.D. N.D.	U	0.2	0.048 0.048	0.024 0.019	N.D. N.D.	U	0.2	0.05	0.023 0.019	NA NA	-	-	-	-	-	-	-	-	-	-
EPA 8270C	Dibenzo[a,h]anthracene	0.0092	0.52	N.D.	Ü	0.2	0.048	0.013	N.D.	U	0.2	0.05	0.013	NA NA	_	-	-	-	-	-	-	-	-	-
	Fluoranthene	1,500	130	N.D.	U	0.2	0.048	0.028	N.D.	U	0.2	0.05	0.027	NA	-	-	-	-	-	-	-	-	-	-
	Fluorene	240	950	0.24		0.2	0.048	0.025	0.23		0.2	0.05	0.024	4.26	-	-	-	-	-	-	-	-	-	-
	Indeno[1,2,3-cd]pyrene	0.092	0.095	N.D.	U	0.2	0.048	0.022	N.D.	U	0.2	0.05	0.022	NA	-	-	-	-	-	-	-	-	-	-
	1-Methylnaphthalene 2-Methylnaphthalene	4.7	10 10	25 20		0.2	0.96 0.96	0.029 0.027	26 22		0.2	0.05	0.028 0.026	3.92 9.52	-	-	-	-	-	-	-	-	-	-
	Naphthalene	17	21	71		2.0	0.96	0.027	76		2.0	0.03	0.020	6.80	-	-	-	-	-	-	-	-	-	-
	Phenanthrene	240	410	N.D.	U	0.2	0.048	0.031	N.D.	U	0.2	0.05	0.030	NA	-	-	-	-	-	-	-	-	-	-
	Pyrene	180	68	N.D.	U	0.2	0.048	0.025	N.D.	U	0.2	0.05	0.025	NA	-	-	-	-	-	-	-	-	-	-
	1,1,1,2-Tetrachloroethane	0.52	50,000	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	NA	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4
	1,1,2,2-Tetrachloroethane 1,1,1-Trichloroethane	0.067 200	500 970	N.D. N.D.	U	1 5	0.5	0.41	N.D. N.D.	U	5	0.5 0.5	0.41	NA NA	N.D. N.D.	U	5	0.5 0.5	0.41	N.D. N.D.	U	1 5	0.5 0.5	0.41
	1,1,2-Trichloroethane	5	50.000	N.D.	U	1	0.5 0.5	0.38	N.D.	U	1	0.5	0.38	NA NA	N.D.	U	1	0.5	0.38	N.D.	U	1	0.5	0.38
	1,1-Dichloroethane	2.4	50,000	N.D.	Ü	5	0.5	0.28	N.D.	Ü	5	0.5	0.28	NA	N.D.	Ü	5	0.5	0.28	N.D.	Ü	5	0.5	0.28
	1,1-Dichloroethylene	7	1,500	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43	NA	N.D.	U	1	0.5	0.43	N.D.	U	1	0.5	0.43
	1,2,3-Trichloropropane	0.6	50,000	N.D.	U	5	1	0.64	N.D.	U	5	1	0.64	NA	N.D.	U	5	1	0.64	N.D.	U	5	1	0.64
	1,2,4-Trichlorobenzene	70	3,000	N.D.	U	5	1	0.5	N.D.	U	5	1	0.5	NA NA	N.D.	U	5	1	0.5	N.D.	U	5	1	0.5
	1,2-Dibromo-3- chloropropane 1,2-Dibromoethane	0.04 0.04	10 50.000	N.D. N.D.	U	10 1	2 0.5	1.2 0.24	N.D. N.D.	U	10	2 0.5	1.2 0.24	NA NA	N.D. N.D.	U	10	0.5	1.2 0.24	N.D. N.D.	U	10	2	1.2 0.24
	1,2-Dibromoetriane 1,2-Dichlorobenzene	600	10	N.D.	U	1	0.5	0.46	N.D.	U	1	0.5	0.46	NA NA	N.D.	U	1	0.5	0.46	N.D.	U	1	0.5 0.5	0.46
	1,2-Dichloroethane	0.15	7,000	N.D.	U	1	0.5	0.24	N.D.	Ü	1	0.5	0.24	NA	N.D.	Ü	1	0.5	0.24	N.D.	Ü	1	0.5	0.24
	1,2-Dichloropropane	5	10	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42	NA	N.D.	U	5	0.5	0.42	N.D.	U	5	0.5	0.42
	1,3-Dichlorobenzene	180	5	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4	NA	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.4
	1,3-Dichloropropene (total of cis/trans)	0.43	50,000	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	NA NA	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25
	1,4-Dichlorobenzene Acetone	75 22,000	5 20,000	N.D. N.D.	U,ICH	20	0.5 10	0.43 6	N.D. N.D.	U,IC	20	0.5 10	0.43 6	NA NA	N.D. 6.5	U J	20	0.5 10	0.43 6	N.D. N.D.	U,IC	20	0.5 10	0.43 6
	Benzene	5	170	N.D.	U	1	0.5	0.14	N.D.	U	1	0.5	0.14	NA NA	N.D.	U	1	0.5	0.14	N.D.	U	1	0.5	0.14
	Bromodichloromethane	0.12	50,000	N.D.	Ü	5	0.5	0.21	N.D.	Ü	5	0.5	0.21	NA	N.D.	Ü	5	0.5	0.21	N.D.	Ü	5	0.5	0.21
	Bromoform	80	510	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5	NA	N.D.	U	10	1	0.5	N.D.	U	10	1	0.5
	Bromomethane	8.7	50,000	N.D.	U	20	5	3.9	N.D.	U,ICJ	20	5	3.9	NA	N.D.	U	20	5	3.9	N.D.	U,ICJ	20	5	3.9
EPA 8260B	Carbon Tetrachloride	5 100	520	N.D.	U	1	0.5	0.23	N.D.	U	1	0.5	0.23	NA NA	N.D.	U	5	0.5	0.23	N.D.	U	1	0.5	0.23
	Chlorobenzene Chloroethane	21,000	50 16	N.D.	U	5 10	0.5 5	0.17 2.3	N.D. N.D.	U	10	0.5 5	0.17 2.3	NA NA	N.D. N.D.	U	10	0.5 5	0.17 2.3	N.D.	U	10	0.5 5	0.17 2.3
	Chloroform	70	2,400	N.D.	Ü	5	0.5	0.46	N.D.	U	5	0.5	0.46	NA NA	N.D.	Ü	5	0.5	0.46	N.D.	U	5	0.5	0.46
	Chloromethane	1.8	50,000	N.D.	U, IJ	10	2	1.8	N.D.	U	10	2	1.8	NA	N.D.	U	10	2	1.8	N.D.	U	10	2	1.8
	cis-1,2-Dichloroethylene	70	50,000	N.D.	U	1	0.5	0.48	N.D.	U	1	0.5	0.48	NA	N.D.	U	1	0.5	0.48	N.D.	U	1	0.5	0.48
	Dibromochloromethane	0.16	50,000	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25	NA NA	N.D.	U	1	0.5	0.25	N.D.	U	1	0.5	0.25
	Ethylbenzene Hexachlorobutadiene	700 0.86	30 6	N.D. N.D.	U	1	0.5 0.5	0.14 0.32	N.D. N.D.	U	1	0.5 0.5	0.14 0.32	NA NA	N.D. N.D.	U	1	0.5 0.5	0.14 0.32	N.D. N.D.	U	1	0.5 0.5	0.14 0.32
	Methyl ethyl ketone (2-Butanone)	7,100	8,400	N.D.	Ü	10	5.0	2.2	N.D.	U	10	5.0	2.2	NA NA	N.D.	Ü	10	5.0	2.2	N.D.	U	10	5.0	2.2
	Methyl isobutyl ketone (4-Methyl-2-Pentanone)	2,000	1300	N.D.	Ü	10	5.0	4.4	N.D.	U	10	5.0	4.4	NA	N.D.	Ü	10	5.0	4.4	N.D.	U	10	5.0	4.4
	Methyl tert-butyl Ether	12	5	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31	NA	N.D.	U	1	0.5	0.31	N.D.	U	1	0.5	0.31
	Methylene chloride	4.8	9,100	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64	NA NA	N.D.	U	5	1.0	0.64	N.D.	U	5	1.0	0.64
ĺ	Styrene 1,1,1,2-Tetrachloroethane	100 0.52	10 50.000	N.D.	U	1	0.5 0.5	0.17 0.4	N.D. N.D.	U	1	0.5 0.5	0.17 0.4	NA NA	N.D. N.D.	U	1	0.5 0.5	0.17 0.4	N.D.	U	1	0.5 0.5	0.17
	1,1,2,2-Tetrachloroethane	0.067	500	N.D.	U	1	0.5	0.41	N.D.	U	1	0.5	0.41	NA NA	N.D.	U	1	0.5	0.4	N.D.	U	1	0.5	0.41
	Tetrachloroethylene	5	170	N.D.	Ü	5	0.5	0.39	N.D.	Ü	5	0.5	0.39	NA	N.D.	U	5	0.5	0.39	N.D.	U	5	0.5	0.39
	Toluene	1,000	40	N.D.	U	1	0.5	0.24	N.D.	U		0.5	0.24	NA	N.D.	U	1	0.5	0.24	N.D.	U	1	0.5	0.24
ĺ	trans-1,2- Dichloroethylene	100	260	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	NA	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37
ĺ	Trichloroethylene	5	310	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37	NA NA	N.D.	U	1	0.5	0.37	N.D.	U	1	0.5	0.37
	Vinyl chloride  Xylenes	10,000	3,400 20	N.D. 0.36	J	1	0.5 0.5	0.3 0.23	N.D. 0.33	J	1	0.5 0.5	0.3 0.23	NA 8.70	N.D. N.D.	U	1 11	0.5 1.5	0.3	N.D. N.D.	U	1 11	0.5 1.5	0.3
EPA 6020	Lead	10,000	50,000	0.36 N.D.	U	1	0.5	0.23	0.33	J	1	0.5	0.23	200	N.D. -	-	-	1.5	-	N.D. -	-	- 11	-	-
	picrograms per liter (ug/L) Shaded values evceeder	•	55,000	11.0.		<u>'</u>	U.L	0.0000	0.170	, ,	'	٥.٤	0.0000	200										•

The data are in micrograms per liter ( $\mu g/L$ ). Shaded values exceeded the DOH EALs.

Not Analyzed

DOH Tier 1 Environmental Action Levels for groundwater where groundwater is a current drinking water source and surface water is greater than 150 meters from the site (DOH, Fall 2011).

Detection Limit or Method Detection Limit (MDL)

Environmental Protection Agency

The chromatographic pattern was inconsistent with the profile of the reference fuel standard. DOH EALS
DL
EPA
HD
ICH
ICJ
J

Initial calibration verification recovery is above the control limit for this analyte. Initial calibration verification recovery is below the control limit for this analyte.

Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated.

LOD LOQ NA N.D. Q TPH-g

Limit of Detection
Limit of Quantitation
Both results for duplicate pair were non-detect, no RPD calculations
Not Detected
Qualifiers

Total Petroleum Hydrocarbons as gasoline
Total Petroleum Hydrocarbons as diesel
Undetected at DL and is reported as less than the LOD.

Contract No. N62742-12-D-1853 Contract Task Order 0002

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#### SECTION 4 - SUMMARY, CONCLUSIONS, AND RECOMMENDATIONS

On July 21 and 22, 2014, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01).

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI. A summary of the analytical results is provided below.

- RHMW01 TPH-d (67 μg/L) was the only analyte detected. The TPH-d concentration detected did not exceed the DOH EALs or the SSRBL.
- RHMW02 TPH-d (1,200 and 1,300 μg/L), TPH-g (48 and 49 μg/L), xylenes (0.36 and 0.33 μg/L), acenaphthene (0.52 and 0.50 μg/L), fluorene (0.24 and 0.23 μg/L), 1-methylnaphthalene (25 and 26 μg/L), 2-methylnaphthalene (20 and 22 μg/L), and naphthalene (71 and 76 μg/L) were detected. TPH-d, 1-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EALs for both drinking water toxicity and gross contamination. However, the TPH-d concentrations did not exceed the SSRBL. 2-Methylnaphthalene was detected at concentrations above the DOH EAL for gross contamination.
- RHMW03 TPH-d (37 µg/L) was the only analyte detected. None of the chemical constituents analyzed for were detected at concentrations above the DOH EALs or the SSRBLs.
- RHMW05 None of the chemical constituents analyzed for were detected.
- RHMW2254-01 None of the chemical constituents analyzed for were detected.

Acetone (6.5 µg/L) was detected in the trip blank submitted on July 21, 2014. Acetone was not detected in any of the groundwater samples, and this is not likely to affect data usability.

#### **Groundwater Contaminant Trends**

RHMW01 – COPCs detected during this round of quarterly sampling are consistent with the
historical data for RHMW01. TPH-d has historically been detected at concentrations above
the DOH EAL for both drinking water toxicity and gross contamination. TPH-d
concentrations have shown an overall decreasing trend from a high of 1,500 μg/L in
February 2005. The TPH-d concentration was last above the DOH EAL for gross
contamination in a sample collected on April 7, 2014.

- RHMW02 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW02. TPH-g, TPH-d, trichloroethylene, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene have historically been detected at concentrations above the DOH EALs. TPH-d concentrations increased sharply from 2,400 µg/L to 5,100 µg/L during the additional sampling event on January 15, 2014, exceeding the SSRBL of 4,500 µg/L. During this July 2014 event, TPH-d concentrations decreased to an average of but remained above both DOH EALs. The concentrations 1,250 µg/L 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene increased to their highest levels since January 2013. The concentrations of TPH-g remained below the DOH EALs for gross contamination and drinking water toxicity. Trichloroethylene has not been detected in RHMW02 since September 2005.
- RHMW03 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW03. TPH-d has historically been detected at concentrations above the DOH EALs; however, it has not been detected at concentrations above the DOH EALs since October 2010.
- RHMW05 No COPCs were detected during this round of quarterly sampling. TPH-d has
  historically been detected in RHMW05 at concentrations above the DOH EAL for both
  drinking water toxicity and gross contamination; however, it has not been detected at
  concentrations above the DOH EALs since January 2010.
- RHMW2254-01 No COPCs were detected during this round of quarterly sampling. TPH-d
  was last detected in RHMW2254-01 at a concentration above the DOH EAL for gross
  contamination in January 2008. Although the method reporting limits for TPH-d exceeded
  one or both DOH EALs between May 2009 and July 2010, TPH-d has not been detected at
  concentrations above the DOH EALs since January 2008.

#### **Conclusions and Recommendations**

Between January and July 2014, the sampling frequency of the wells and monitoring point inside Red Hill increased, due to response activities related to a reported release at Tank 5. Since the previous quarterly sampling event in April 2014, three wells (RHMW01, RHMW02, and RHMW05) and one sampling point (RHMW2254-01) have each been sampled three times, including this quarterly event. During this period only TPH-d in RHMW01 and TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 were detected at concentrations exceeding the DOH EALs. However, the TPH-d concentration detected in RHMW01 only exceeded the DOH EAL during the April 7, 2014 release response groundwater monitoring event. Groundwater contaminant concentrations in the other three wells (RHMW03, RHMW05, and RHMW2254-01) remained at low concentrations and did not change significantly, or were not detected.

Concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 show an increasing trend since March 5, 2014, but remain consistent with historical data. All other analytical results were generally consistent with historical data.

Based on the groundwater monitoring results and the reported release at Tank 5 in January 2014, continued groundwater monitoring at the wells inside the RHSF tunnel is recommended. The next quarterly event is tentatively scheduled for October 2014.

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#### **SECTION 5 – FUTURE WORK**

Future work includes the fourth quarter 2014 groundwater monitoring which is tentatively scheduled for October 2014. A quarterly groundwater monitoring report will be prepared to document the sampling event.

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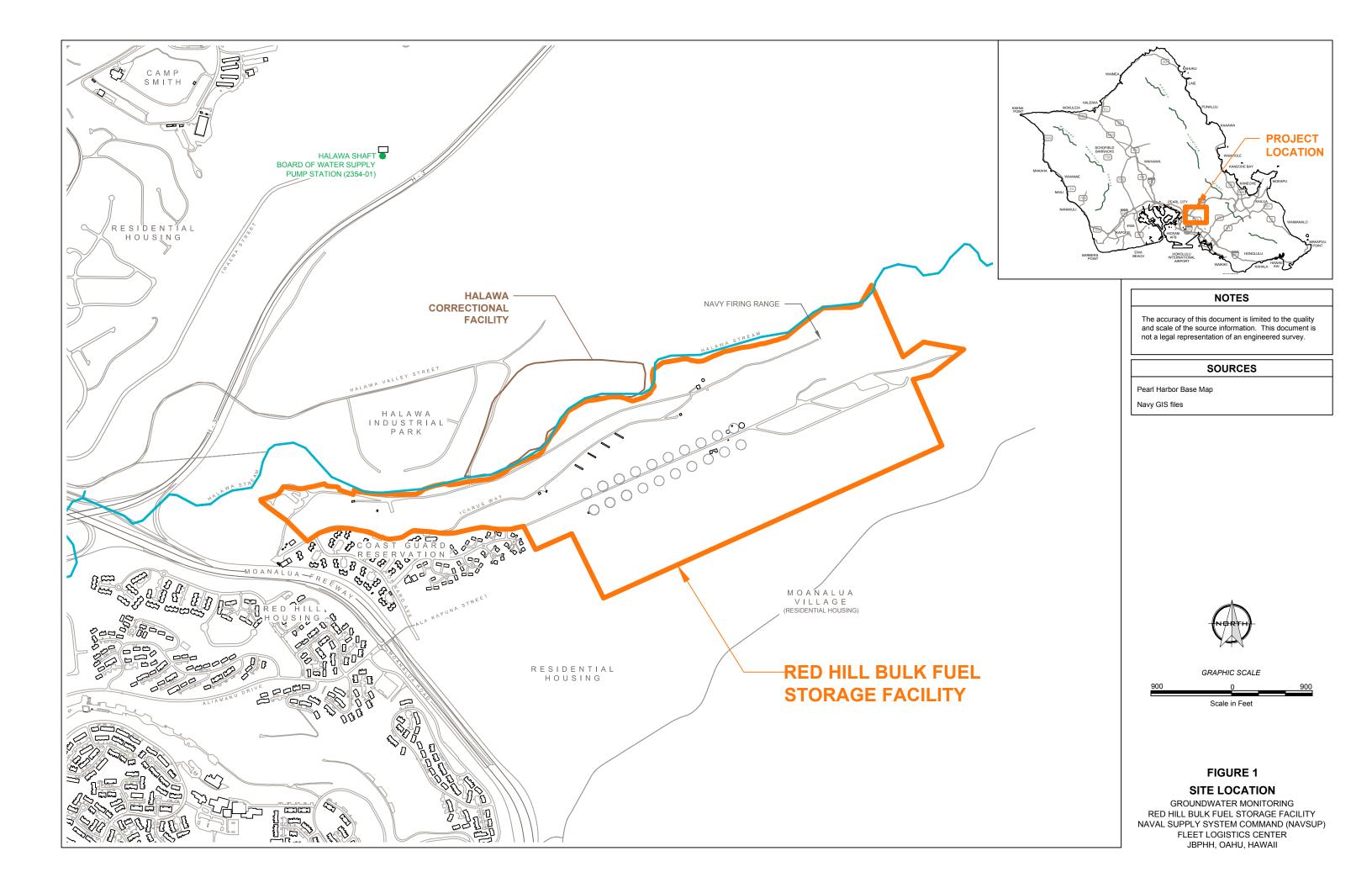
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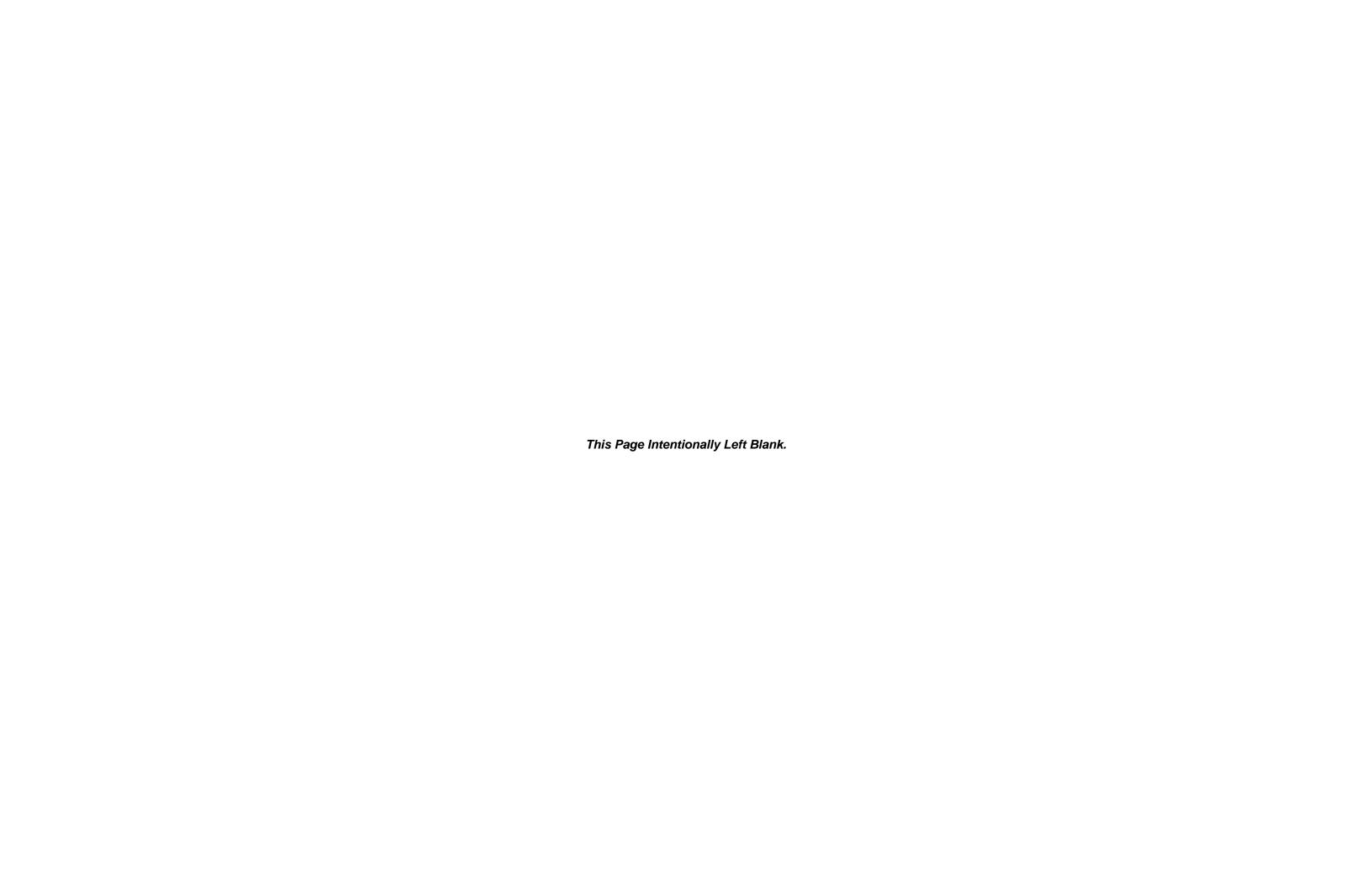
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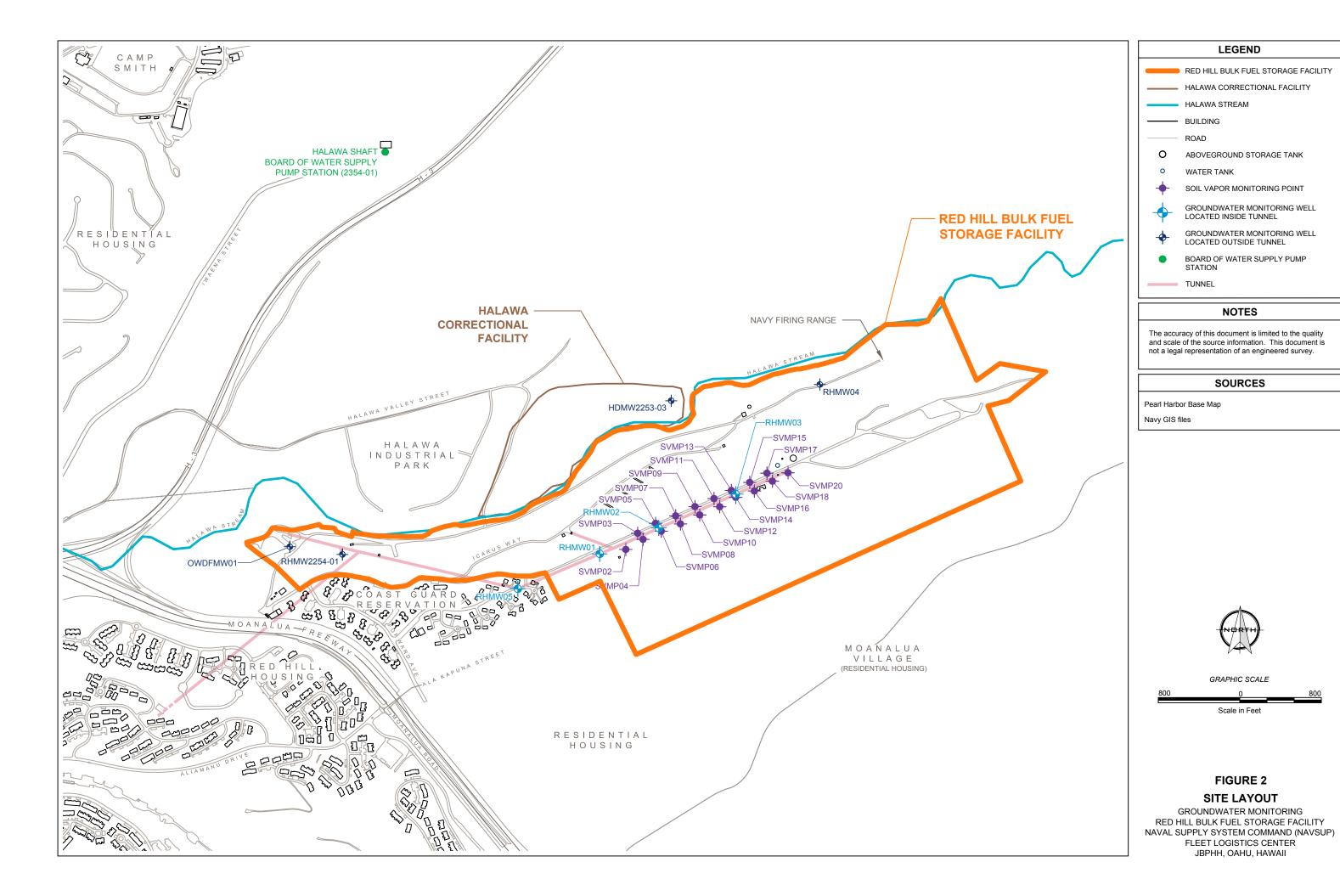
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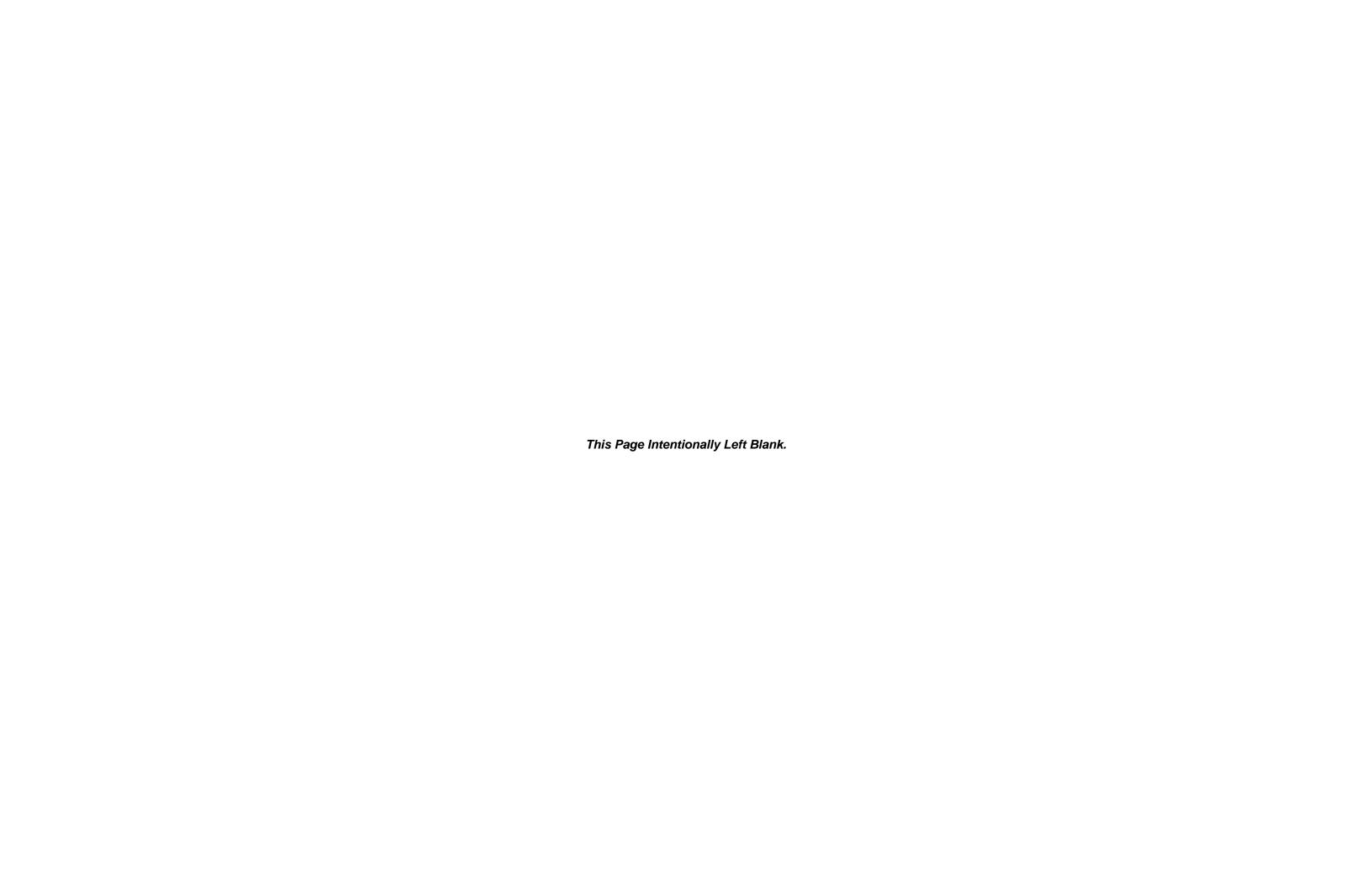
## **FIGURES**











# APPENDIX A Groundwater Sampling Logs





Well ID:	RHMW01	Locat	ion: Red Hill	Bulk Fuel Stora	age Facility P	roject No.:	112066	
Initial W	ater Level:	84.13 ft	Date:	7/21/2014	т	ime: 830		
Total De	epth of Well:	97.35 ft	Person	nel Involved:	Justin Lan	n, Jeff Hatteme	<u>r</u>	
Length o	of Saturated Z	one: 13.22	ft Weathe	Weather Conditions: Not applicab		ole – well is located indoors		
Volume	of Water to be	e Removed: 2	.5 L Method	of Removal:	Blad	der Pump		
Water L	evel After Pur	ging: <u>84.13</u>	ft Pumpin	ng Rate:	0.17	L/min		
Well Pu	rge Data:							
	Volum		Conductivity				Redox (ORP)	
Time	Remove	ed pH	(mS/cm)	DO (mg/l)	Temperature	Salinity	(mV)	
837	0.0 L	7.77	0.343	7.60	25.6		95.1	
840	0.5 L	7.20	0.35	7.12	25.33		56.1	
843	1.0 L	7.12	0.351	4.07	25.18	-	36.2	
846	1.5 L	7.07	0.352	2.33	25.00	-	-6.7	
849	2.0 L	7.05	0.352	2.3	25.05	-	-7.1	
854	2.5 L	7.06	0.352	2.28	25.10		-7.3	
Sample	Withdrawal M	lethod:	Bladder Pur	mp				
Appeara	ince of Sampl	e:						
	Color:		Clear					
	Turbidity	<i>r</i> :	Low					
	Sedimer	nt:	None					
	Other:		None		•			
Laborato	ory Analysis P	arameters and F	Preservatives:		; TPH-g, VOCs - 82	260; PAHs - 82	70c sim;	
				lead - 6020				
Number	and Types of	Sample Contain	ers: <u>6 - 40m</u>	nl VOAs, 2 - 1L a	amber jar, 1 - 500m	nl amber jar, 1	- 250ml HDPE	
Sample	Identification	Numbers: ES	3103 [0855]				_	
Deconta	mination Prod	cedures: Triple	Rinsed					
Notes:	YSI did not h	ave salinity para	meter.					
Sample		ı Lam, Jeff Hatte						
•	d Delivered to	: Calscien	ce Environmenta	al Lab	Transporters: Fed	Ex		
Date:	7/21/2014		0 "	-1 (0 !!	Time: 1200			
			Capacity of Cas	sing (Gallons/Li	near ⊦eet)			



Well ID: R	HMW02	Location:	Red Hill	Bulk Fuel Stora	ge Facility P	roject No.:	112066
Initial Water	Level: 86	80 ft	Date:	7/21/2014	т	ime: 936	
Total Depth	of Well:	92.91 ft	Personn	nel Involved:	Justin Lam	n, Jeff Hattemer	
Length of Sa	aturated Zone:	6.11 ft	Weathe	r Conditions:	Not applicable –	well is located	indoors
Volume of V	Vater to be Re	moved: <u>2.0 L</u>	Method	of Removal:	Blad	der Pump	
Water Leve	After Purging	86.80 ft	_ Pumping	g Rate:	0.17	L/min	
Well Purge							
Time	Volume Removed		onductivity (mS/cm)	DO (mg/l)	Temperature	Salinity	Redox (ORP) (mV)
944	0.0 L	7.09	0.581	0.50	24.12	<u> </u>	-163.9
947	0.5 L	7.07	0.582	0.48	24.04		-162.7
950	1.0 L	7.05	0.585	0.41	24.00		-151.8
953	1.5 L	7.04	0.590	0.40	23.95	-	-150.6
953	2.0 L	7.04	0.590	0.40	23.96		-149.7
Sample Wit	hdrawal Metho	od:	Bladder Pum	ıp			
Appearance	of Sample:			_			
	Color:		Clear				
	Turbidity:		Low				
	Sediment:		None	_			
	Other:		None				
Laboratory /	Analysis Paran	neters and Prese	rvatives:	TPH-d - 8015;	TPH-g, VOCs - 82	260; PAHs - 827	0c sim;
				lead - 6020			
Number and	d Types of San	nple Containers:	16 - 40ml V	/OAs, 6 - 1L am	ber jar, 4 - 500ml	amber jar, 4 - 50	00ml HDPE
Sample Idea	ntification Num	bers: ES104	[1000], ES10	04 MS/MSD [10	00], ES105 (Dup)	[1055]	
Decontamin	ation Procedu	res: Triple Rins	sed				
		salinity paramete	r.				
Sampled by		n, Jeff Hattemer		11 -1-	F	=	
Sampled De	_	Calscience E	nvironmenta		Transporters: FedI	±Χ	
Date: <u>7/2</u>	1/2014	Car	pacity of Cas	ing (Gallons/Lin	Time: <u>1200</u> ear Feet)		



Well ID: R	HMW03	Location:	Red Hill	Bulk Fuel Stora	age Facility	Project No.:	112066
Initial Water	Level: 10	02.98 ft	Date:	7/22/2014		Time: 11	100
Total Depth	of Well:	110.12 ft	Personr	nel Involved:	Justin I	₋am, Jeff Hatter	ner
Length of Sa	aturated Zor	ne: 7.14 ft	Weathe	er Conditions:	Not applicab	le – well is locat	ed indoors
Volume of V	Vater to be F	Removed: 4.0 L	Method	of Removal:	В	ladder Pump	
Water Leve	l After Purgir	ng: <u>102.98 ft</u>	Pumpin	g Rate:	C	0.33 L/min	
Well Purge							
Time	Volume Removed		onductivity (mS/cm)	DO (mg/l)	Temperature	e Salinity	Redox (ORP) (mV)
1105	0.0 L	6.71	0.743	2.40	26.90		123.5
1108	1.0 L	6.72	0.742	2.25	26.87		110.5
1111	2.0 L	6.76	0.746	1.64	26.61		84.0
1114	3.0 L	6.76	0.746	1.60	26.64		79.9
1117	4.0 L	6.78	0.746	1.61	26.60		76.0
						<u> </u>	
						<u> </u>	
						_	
						_	
Sample Wit	hdrawal Met	hod:	Bladder Pur	np			
Appearance	of Sample:						
	Color:		Clear				
	Turbidity:		Low				
	Sediment:		None				
	Other:		None				
Laboratory A	Analysis Par	ameters and Prese	ervatives:	TPH-d - 8015;	TPH-g, VOCs	- 8260; PAHs - 8	8270c sim;
·	•			lead - 6020	<u> </u>		·
Number and	d Types of S	ample Containers:	6 - 40m	I VOAs, 2 - 1L a	amber jar, 1 - 50	00ml amber jar,	1 - 250ml HDPE
Sample Idea	ntification Nu	umbers: ES106	[1120]				
Decontamin	ation Proced	dures: Triple Rins	sed				
Notes: YS	I did not hav	e salinity paramete	er.				
Sampled by	: Justin L	.am, Jeff Hattemer					
Sampled De		Calscience E	nvironmenta		Transporters: F	edEx	
Date: <u>7/2</u>	2/2014	Cor	noity of Co	sing (Gallons/Li	Time: 1215		
		Car	Jaciiv Ol Cas	sina walions/Lli	пеаг гееп		



Well ID: F	RHMW05	Location:	Red Hill	Bulk Fuel Stora	age Facility	Project No.:	112066
Initial Wate	r Level:	83.49 ft	Date:	7/22/2014		Time: 1013	3
Total Depth	of Well:	Unable to Measure	Personr	nel Involved:	Justin La	am, Jeff Hatteme	<u>r</u>
Length of S	aturated Zo	ne: Unknown	_ Weathe	er Conditions:	Not applicable	e – well is located	indoors
Volume of \	Water to be	Removed: 4.0 L	Method	of Removal:	Bla	adder Pump	
Water Leve	l After Purgi	ing: 83.49 ft	Pumpin	g Rate:	0.5	33 L/min	
Well Purge	Data:						
	Volume		onductivity	<b></b>		<b>.</b>	Redox (ORP)
Time	Removed		(mS/cm)	DO (mg/l)	Temperature	Salinity	(mV)
1018	0.0 L	7.73	0.837	9.60	23.67	. <u> </u>	98.9
1021	1.0 L	<u>7.64</u>	0.839	8.42	23.40	. <del></del>	90.7
1024	2.0 L	7.58	0.836	7.98	22.79	<u> </u>	68.9
1027	3.0 L	7.54	0.836	7.80	22.68	<u> </u>	67.1
1030	4.0 L	7.53	0.830	7.80	22.66	<u> </u>	66.8
						<u> </u>	<u> </u>
						<u> </u>	_
					-		
Sample Wit	thdrawal Me	thod:	Bladder Pur	mp			
Appearance	e of Sample						
	Color:		Clear				
	Turbidity:		Clear				
	Sediment		None	,			
	Other:		None				
				_			
Laboratory	Analysis Pa	rameters and Prese	ervatives:	TPH-d - 8015;	; TPH-g, VOCs -	8260; PAHs - 82	70c sim;
				lead - 6020			
Number and	d Types of S	Sample Containers:	6 - 40m	I VOAs, 2 - 1L a	amber jar, 1 - 500	oml amber jar, 1	- 250ml HDPE
Sample Ide	ntification N	umbers: ES108	3 [1035]				
Decontamir	nation Proce	dures: Triple Rin	sed				
Notes: YS	I did not hav	ve salinity paramete	er.				
Sampled by		Lam, Jeff Hattemer					
Sampled D		Calscience E			Transporters: Fe	edEx	
Date: <u>7/2</u>	22/2014				Time: 1215		
		Ca <sub>l</sub>	pacity of Cas	sing (Gallons/Li	near Feet)		



Well ID	: RHMW2254	<u>-01</u> Loca	tion: Red H	ill Bulk Fuel Stor	age Facility F	Project No.:	112066
Initial W	/ater Level:	82.19 ft	Date:	7/22/2014		ime: 915	
Total D	epth of Well:	Not applica	ble Persor	nnel Involved:	Justin Lar	n, Jeff Hatteme	<u>r</u>
Length	of Saturated Z	one: Not appli	cable Weath	er Conditions:	Not applicable -	- well is located	indoors
Volume	of Water to b	e Removed: 4	.0 L Metho	d of Removal:	Blac	lder Pump	
Water L	evel After Pur	ging: 82.	19 ft Pumpi	ng Rate:	0.33	3 L/min	
Well Pu	ırge Data:						
	Volum		Conductivity				Redox (ORP)
Time		·	(mS/cm)	DO (mg/l)	Temperature	Salinity	(mV)
920			0.504	7.92	23.03		45.4
923	1.0 L	7.34	0.525	7.55	22.52	_	29.1
926	2.0 L	7.30	0.527	7.53	22.44		27.9
929	3.0 L	7.32	0.527	7.52	22.41		27.7
932	4.0 L	7.31	0.527	7.52	22.41		27.6
			- ·	·			
			_	<del></del>			
Sample	Withdrawal M	lethod:	Bladder Pu	ımp			
Appear	ance of Sampl	e:					
	Color:		Clear				
	Turbidity	/:	Clear				
	Sedimer	nt:	None		•		
	Other:		None		•		
Laborat	ory Analysis P	arameters and	Preservatives:	TPH-d - 8015	; TPH-g, VOCs - 8	260; PAHs - 82 <sup>°</sup>	70c sim;
				lead - 200.8			
Numbe	r and Types of	Sample Contai	ners: 6 - 40r	ml VOAs, 2 - 1L	amber jar, 1 - 500n	nl amber jar, 1 -	250ml HDPE
Sample	Identification	Numbers: E	S107, ES107 U	F [0945]			
Deconta	amination Prod	cedures: Triple	e Rinsed				
Notes:	YSI did not h	ave salinity para	ameter.				
Sample		n Lam, Jeff Hatt					
•	d Delivered to	: Calscier	nce Environmen	tal Lab	Transporters: Fed	Ex	-
Date:	7/22/2014				Time: 1215		
			Capacity of Ca	asing (Gallons/Li	near Feet)		



# APPENDIX B Field Notes



58 R175T Project / Client NAVFAC Personnel JIJ Softy not 0800 @ RHST- Calibrate 1102 Softy noty 0810 Byler June 1 0830 @ RHMWOI, DTW = 84,13' b foc. 0855 Collected \$5103, 2,52,000,00 0936 @ RHMWOZ, DTW 86,80 How. and duplicated ES105 (recorded) 7000 as 1955). .

Location R	135-	Date 7/2/14	59
Project / Client _	NAVEAL	30.0	

			1 1		0 -	
	082	Begin	n purge	SVA	1605	4
	Shallow	556	513	481	1556	527
	Mid	423	382	339	423	1392
	Deep	426	397	440	1440	426
	0834	Begin	purge	SUM	803	
	Shaller	922	1000	1078	1084	1021
	mid	1032	1162	1133	1203	1133
	deep	1136	1200	1261	1261	1215
	0856	Begin	Quig	e 50,	nroy.	
	shallor	571	522	\$65	1571	1557
	mid	620	556	626	698	625
	desp	1797	707	63C	797	735
	0914	Begin	proje	SUMJO	+	
	Shiller	107	- 11)	113	113	1 1
P	len of	237	232	231	237	234
<del>)</del>	der	239	135	233	239	237
	0941	Begin	mose	SUN	1706	
	Shallor	20.6	219	22,5	122,5	121.91
4	mid 1	20,1	21.6	223	1223	121.61
	0958	Begin	ywa	SVI	RO7	
y	Shilly	33,0	35,41	35,5	135.5	134.9
	mid)	Sherkeel.	No	Sample		
A	deep	23.8	25,3	25,3	253	1249
	A	c 88m				

Location RHSP Date 7/21/11 Project / Client NA VZAC SUMPOS 10/08 Begin purge 13.7 13.7 13.2 16.9 16.9 16.5 17.8 17.8 16.8 SUMP 09 Shellor 13,2 123 16.5 15.7 My 17,0 14,6 Deep Begin purce 1035 5957 | 6450 | 6714 | 6714 | 6459 Shaller 10,2 10,2 9981 4652 9872 nd 12.6 | 12.0 10,8 12.0 12.6 der SIMPID .. Blan yunge 1051 Shallow 3123 3346 3433 3433 3347 day 3494 3622 3722 3722 3640 1120 Deport Time!.
1150 Pack camples, dump of &
water : e Adit 3 had to d aff samples @ fel by 5.00 propord 1225

Location R17 5 F. Project / Client MXVXX Z	Date 7/22/14 61
Project / Client	
1 Toject / Oliciti	

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Var	2 324	NO	1 4		11	. )		_							
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		5	nfe	5		m	ee)	~~	>			1	19-1		
77	SE	)	13	nte	<b>Y</b>	1	wr	ne	1.						
08	30		<b>SC</b>	.P	ורעוט	9	1	ימס	, e				0		lad.
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	14	N										1			
09	15	0		RH	Mh	122	5	1-1	11,	07	V =	87	-110	1, 9	166.
179	45		Co	110	+-1	9	90	my	C	E	510	7	E.	107	U
	13	6	2 1	2 h .	MW	105	5	D	W	= 9	33	49	3	to	19 5
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Project / Client N&VY-N C

18 N.V.

					and the second second
0805	Begin	prige	5	SMR20	
Shellow	1454	1 1551	116/1	1 167)	1158/
mid	1751	1906	1889	1960	1877
Leon	2057	2274	12246	1960	2213
0815	Benin	purge	311	7818	
Shallow	1243	1423	11614	1614	11474
deep	12454	2812	3012	13012	12823
0842	Begin			MPIZ	
Shallon	1346	1586	1689	1689	11578
mid	1823	2103	2349	2349	12,156
deg	2403	2760	3026	3026	2804
0904	Begn	frese	5	1 M9 16	
shallow 1	477	560	1585	1 585	552
mid	391	440	494	1494	455
deg	1414	1 431	554	554	488
0924	Begn	pure	x 5v	MP 15	
Shallor	6111	6341	703	MP 15	663
m'd	mid	block	d. No	somale	-
deep	11151	1306	1414	1414	1321
0938	polyn	proje	5	MP 14	
Shallow	277	305	351	360	323
md	374	482			LATY
Jest 1	531	594			634

Location RHSF
Project / Client NAVFAC

0955 proje SVMP13 Begin 531 505 | 531 Shallon 495 414 mid 637 800 911 1103 1211 11225 11225 deep 1013 Smllw Begin 51MP 12 purge 242 | 237 | 290 | 222 | 194 | 228 302 | 311 | 317 160 2421 mid 290 deep 1030 Ben prize SVMPII 225 | 242 | 231 mid 1261 240 blocked No deep 1145 Depart funne) packing samples, ドinish water dump ous, drum @ 1215 Depart Fed Ex to



# APPENDIX C Laboratory Reports





### Calscience



## **WORK ORDER NUMBER: 14-07-1466**

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

**Analytical Report For** 

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

**Attention:** Domonkos Feher, Ph.D.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Richard Vellas

Approved for release on 07/30/2014 by:

Richard Villafania Project Manager



ResultLink >

Email your PM >

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.



### **Contents**

Client Project Name: Red Hill LTM 112066 Work Order Number: 14-07-1466

1	Work Order Narrative	3
2	Client Sample Data.  2.1 EPA 8015B (M) TPH Diesel (Aqueous).  2.2 EPA 6020 ICP/MS Metals (Aqueous).  2.3 EPA 8270C SIM PAHs (Aqueous).  2.4 GC/MS GRO/EPA 8260B Volatile Organics (Aqueous).	4 4 5 6 10
3	Quality Control Sample Data.  3.1 MS/MSD.  3.2 PDS/PDSD.  3.3 LCS/LCSD.	20 20 25 26
4	Sample Analysis Summary	31
5	Glossary of Terms and Qualifiers	32
6	Chain-of-Custody/Sample Receipt Form	33



#### **Work Order Narrative**

Work Order: 14-07-1466 Page 1 of 1

#### **Condition Upon Receipt:**

Samples were received under Chain-of-Custody (COC) on 07/22/14. They were assigned to Work Order 14-07-1466.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

#### **Holding Times:**

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

#### **Quality Control:**

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

#### **Additional Comments:**

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New\_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

#### **Subcontractor Information:**

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Page 1 of 1



Project: Red Hill LTM 112066

#### **Analytical Report**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1466

Kailua, HI 96734-2500

Preparation:

Method:

Units:

Units:

Date Received:

07/22/14

Work Order:

14-07-1466

EPA 3510C

EPA 8015B (M)

Units:

ug/L

Client Sample Number		Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES103		14-07-1466-1-I	07/21/14 08:55	Aqueous	GC 45	07/23/14	07/24/14 21:13	140723B11A
Comment(s):	- Results were evaluated to	the MDL (DL), con	centrations >= to t	the MDL (DL	_) but < RL (LO	Q), if found, ar	e qualified with a	a "J" flag.
<u>Parameter</u>		Result	<u>DL</u>	<u>LOD</u>	LOC	<u>2</u>	<u>DF</u>	<b>Qualifiers</b>
TPH as Diesel		67	11	12	25		1.00	HD
Surrogate		Rec. (%)	Control Limi	<u>its Qualit</u>	fiers			
n-Octacosane		127	51-141					

ES104	14-07-1466-2-	l 07/21/14 10:00	Aqueous	GC 45		/24/14 140723B11 :30	A
Comment(s):	- Results were evaluated to the MDL (DL),	concentrations	>= to the MDL (D	L) but < RL (LOQ)	, if found, are qual	fied with a "J" flag.	
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<b>Qualifiers</b>	
TPH as Diesel	1200	11	12	25	1.00	HD	
<u>Surrogate</u>	<u>Rec. (%</u>	<u>(6)</u> <u>Cont</u>	rol Limits Qual	<u>ifiers</u>			
n-Octacosane	110	51-14	41				

ES105	14-07-		07/21/14 / 10:55	Aqueous	GC 45	07/23/14	07/24/14 21:48	140723B11A
Comment(s):	- Results were evaluated to the MD	L (DL), concer	ntrations >= to th	ne MDL (DL)	) but < RL	(LOQ), if found, a	re qualified with a	"J" flag.
<u>Parameter</u>		Result	<u>DL</u>	<u>LOD</u>		<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
TPH as Diesel		1300	11	12		25	1.00	HD
Surrogate n-Octacosane		Rec. (%) 135	Control Limit 51-141	ts Qualifi	<u>ers</u>			

Method Blank	099-15-516-158	N/A Aq	ueous GC 45	5 07/23/14	07/24/14 18:11	140723B11A
Comment(s): - Results were ev	valuated to the MDL (DL), conc	entrations >= to the I	MDL (DL) but <	RL (LOQ), if found,	are qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
TPH as Diesel	<12	11	12	25	1.00	U
<u>Surrogate</u>	Rec. (%)	Control Limits	<u>Qualifiers</u>			
n-Octacosane	101	51-141				



Method Blank

#### **Analytical Report**

Environmental Science International, Inc. Date Received: 07/22/14 354 Uluniu Street, Suite 304 Work Order: 14-07-1466 Preparation: EPA 3005A Filt. Kailua, HI 96734-2500 Method: EPA 6020 Units: mg/L

Page 1 of 1 Project: Red Hill LTM 112066

Client Sample I	Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES103		14-07-1466-1-J	07/21/14 08:55	Aqueous	ICP/MS 04	07/23/14	07/24/14 19:50	140723L03D
Comment(s):	- Results were evaluated t	o the MDL (DL), cond	centrations >= to	the MDL (DL	) but < RL (LC	Q), if found, a	re qualified with a	a "J" flag.
<u>Parameter</u>		Result	<u>DL</u>	<u>LOD</u>	LC	<u>Q</u>	<u>DF</u>	<u>Qualifiers</u>
Lead		< 0.000200	0.0000898	0.000	200 0.0	0100	1.00	U

ES104	14-07-1466-2-J	07/21/14 10:00	Aqueous	ICP/MS 04		07/24/14 140723L03D 19:48
Comment(s):	- Results were evaluated to the MDL (DL), cor	centrations >= to	the MDL (DL	) but < RL (LOC	), if found, are qu	ualified with a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>DF</u>	<u>Qualifiers</u>
Lead	<0.000200	0.0000898	0.0002	0.00	100 1.00	) U

ES105	14-07-1466-3-J	07/21/14 10:55	Aqueous	ICP/MS 04	07/23/14	07/24/14 19:52	140723L03D
Comment(s):	- Results were evaluated to the MDL (DL), con	centrations >= to	the MDL (DL	) but < RL (LO	Q), if found,	are qualified with a	"J" flag.
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOC	<u>2</u>	<u>DF</u>	<u>Qualifiers</u>
Lead	0.000170	0.0000898	0.0002	200 0.00	0100	1.00	J

Aqueous ICP/MS 04

07/23/14

07/24/14

						19:24	
Comment(s):	- Results were evaluated to the N	MDL (DL), concer	ntrations >= to the	MDL (DL) but <	RL (LOQ), if foun	d, are qualified	with a "J" flag.
<u>Parameter</u>		<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
Lead		<0.000200	0.0000898	0.000200	0.00100	1.00	U

099-14-497-87



140723L03D



### **Analytical Report**

Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

07/22/14 14-07-1466

Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 1 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID	
ES103	14-07-1466-1-G	07/21/14 08:55	Aqueous	GC/MS AAA	07/28/14	07/29/14 16:17	140728L01	
Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.								
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	<u>Qualifiers</u>	
Naphthalene	<0.051	0.024	0.051	0.20	)	1.00	U	
2-Methylnaphthalene	<0.051	0.027	0.051	0.20	)	1.00	U	
1-Methylnaphthalene	<0.051	0.029	0.051	0.20	)	1.00	U	
Acenaphthylene	<0.051	0.018	0.051	0.20	)	1.00	U	
Acenaphthene	< 0.051	0.021	0.051	0.20	)	1.00	U	
Fluorene	<0.051	0.025	0.051	0.20	)	1.00	U	
Phenanthrene	<0.051	0.031	0.051	0.20	)	1.00	U	
Anthracene	< 0.051	0.035	0.051	0.20	)	1.00	U	
Fluoranthene	<0.051	0.028	0.051	0.20	)	1.00	U	
Pyrene	< 0.051	0.025	0.051	0.20	)	1.00	U	
Benzo (a) Anthracene	< 0.051	0.024	0.051	0.20	)	1.00	U	
Chrysene	<0.051	0.019	0.051	0.20	)	1.00	U	
Benzo (k) Fluoranthene	< 0.051	0.024	0.051	0.20	)	1.00	U	
Benzo (b) Fluoranthene	< 0.051	0.025	0.051	0.20	)	1.00	U	
Benzo (a) Pyrene	<0.051	0.037	0.051	0.20	)	1.00	U	
Indeno (1,2,3-c,d) Pyrene	<0.051	0.022	0.051	0.20	)	1.00	U	
Dibenz (a,h) Anthracene	< 0.051	0.027	0.051	0.20	)	1.00	U	
Benzo (g,h,i) Perylene	<0.051	0.022	0.051	0.20	)	1.00	U	
<u>Surrogate</u>	Rec. (%)	Control Limi	its Qualifi	<u>iers</u>				
Nitrobenzene-d5	75	28-139						
2-Fluorobiphenyl	76	33-144						
p-Terphenyl-d14	73	23-160						



Environmental Science International, Inc.

354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

Work Order: 14-07-1466

Preparation:

EPA 8270C SIM PAHs

Method: Units:

ug/L

07/22/14

EPA 3510C

Project: Red Hill LTM 112066

Page 2 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES104	14-07-1466-2-G	07/21/14 10:00	Aqueous	GC/MS AAA	07/28/14	07/29/14 16:41	140728L01
Comment(s): - Results were evaluated to	o the MDL (DL), con	centrations >= to t	he MDL (DL	) but < RL (LOC	Q), if found, a	re qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	<u>Qualifiers</u>
Acenaphthylene	<0.048	0.017	0.048	0.19	)	1.00	U
Acenaphthene	0.52	0.020	0.048	0.19	)	1.00	
Fluorene	0.24	0.023	0.048	0.19	)	1.00	
Phenanthrene	<0.048	0.029	0.048	0.19	)	1.00	U
Anthracene	<0.048	0.033	0.048	0.19	)	1.00	U
Fluoranthene	<0.048	0.026	0.048	0.19	)	1.00	U
Pyrene	<0.048	0.024	0.048	0.19	)	1.00	U
Benzo (a) Anthracene	<0.048	0.023	0.048	0.19	)	1.00	U
Chrysene	<0.048	0.018	0.048	0.19	)	1.00	U
Benzo (k) Fluoranthene	<0.048	0.022	0.048	0.19	)	1.00	U
Benzo (b) Fluoranthene	<0.048	0.024	0.048	0.19	)	1.00	U
Benzo (a) Pyrene	<0.048	0.035	0.048	0.19	)	1.00	U
Indeno (1,2,3-c,d) Pyrene	<0.048	0.021	0.048	0.19	)	1.00	U
Dibenz (a,h) Anthracene	<0.048	0.026	0.048	0.19	)	1.00	U
Benzo (g,h,i) Perylene	<0.048	0.021	0.048	0.19	)	1.00	U
Surrogate	Rec. (%)	Control Limi	its Qualifi	<u>iers</u>			
Nitrobenzene-d5	82	28-139					
2-Fluorobiphenyl	79	33-144					
p-Terphenyl-d14	79	23-160					

ES104	14-07-1466-2-G	07/21/14 A 10:00	Aqueous G	C/MS AAA 07/28/1	4 07/29/14 21:31	140728L01
Comment(s): - Results were evalu	ated to the MDL (DL), conc	entrations >= to the	e MDL (DL) b	out < RL (LOQ), if four	nd, are qualified wit	h a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<b>Qualifiers</b>
Naphthalene	71	0.44	0.96	3.8	20.0	
2-Methylnaphthalene	20	0.51	0.96	3.8	20.0	
1-Methylnaphthalene	25	0.54	0.96	3.8	20.0	
Surrogate	Rec. (%)	Control Limits	<u>Qualifier</u>	<u>s</u>		
Nitrobenzene-d5	65	28-139				
2-Fluorobiphenyl	77	33-144				
p-Terphenyl-d14	52	23-160				



Environmental Science International, Inc.

354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

d: 07/22/14 14-07-1466

Preparation:

Work Order:

EPA 8270C SIM PAHs

Method: Units:

ug/L

Offics.

Page 3 of 4

EPA 3510C

Project: Red Hill LTM 112066

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES105	14-07-1466-3-G	07/21/14 10:55	Aqueous	GC/MS AAA	07/28/14	07/29/14 17:05	140728L01
Comment(s): - Results were evaluated t	o the MDL (DL), con	centrations >= to t	the MDL (DL	.) but < RL (LOC	Q), if found, a	are qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOC	<u>2</u>	<u>DF</u>	<u>Qualifiers</u>
Acenaphthylene	<0.051	0.018	0.051	0.20	)	1.00	U
Acenaphthene	0.50	0.021	0.051	0.20	)	1.00	
Fluorene	0.23	0.025	0.051	0.20	)	1.00	
Phenanthrene	<0.051	0.031	0.051	0.20	)	1.00	U
Anthracene	<0.051	0.035	0.051	0.20	)	1.00	U
Fluoranthene	<0.051	0.028	0.051	0.20	)	1.00	U
Pyrene	<0.051	0.025	0.051	0.20	)	1.00	U
Benzo (a) Anthracene	<0.051	0.024	0.051	0.20	)	1.00	U
Chrysene	<0.051	0.019	0.051	0.20	)	1.00	U
Benzo (k) Fluoranthene	<0.051	0.024	0.051	0.20	)	1.00	U
Benzo (b) Fluoranthene	< 0.051	0.025	0.051	0.20	)	1.00	U
Benzo (a) Pyrene	<0.051	0.037	0.051	0.20	)	1.00	U
Indeno (1,2,3-c,d) Pyrene	<0.051	0.022	0.051	0.20	)	1.00	U
Dibenz (a,h) Anthracene	< 0.051	0.027	0.051	0.20	)	1.00	U
Benzo (g,h,i) Perylene	<0.051	0.022	0.051	0.20	)	1.00	U
Surrogate	Rec. (%)	Control Lim	its Qualif	<u>iers</u>			
Nitrobenzene-d5	85	28-139					
2-Fluorobiphenyl	80	33-144					
p-Terphenyl-d14	79	23-160					

ES105	14-07-1466-3-G	07/21/14 Ac 10:55	queous G	C/MS AAA	07/28/14	07/29/14 21:55	140728L01
Comment(s): - Results were evaluated to	o the MDL (DL), cond	centrations >= to the	MDL (DL) b	ut < RL (LOQ	), if found, are	qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOQ	<u> </u>	<u>)F</u>	<b>Qualifiers</b>
Naphthalene	76	0.47	1.0	4.1	2	20.0	
2-Methylnaphthalene	22	0.54	1.0	4.1	2	20.0	
1-Methylnaphthalene	26	0.57	1.0	4.1	2	20.0	
Surrogate	Rec. (%)	Control Limits	Qualifiers	<u>s</u>			
Nitrobenzene-d5	57	28-139					
2-Fluorobiphenyl	65	33-144					
p-Terphenyl-d14	47	23-160					



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

07/22/14 14-07-1466

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank	099-15-148-52	N/A	Aqueous	GC/MS AAA	07/28/14	07/29/14 13:03	140728L01
Comment(s): - Results were evaluated to	the MDL (DL), con	centrations >= to tl	he MDL (DL	) but < RL (LOC	), if found, are	qualified with a	"J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOC	<u>]</u>	<u>DF</u>	<u>Qualifiers</u>
Naphthalene	< 0.050	0.023	0.050	0.20	1	1.00	U
2-Methylnaphthalene	< 0.050	0.026	0.050	0.20	1	1.00	U
1-Methylnaphthalene	< 0.050	0.028	0.050	0.20	1	1.00	U
Acenaphthylene	< 0.050	0.018	0.050	0.20	1	1.00	U
Acenaphthene	< 0.050	0.021	0.050	0.20	1	1.00	U
Fluorene	< 0.050	0.024	0.050	0.20	1	1.00	U
Phenanthrene	< 0.050	0.031	0.050	0.20	1	1.00	U
Anthracene	< 0.050	0.034	0.050	0.20	1	1.00	U
Fluoranthene	< 0.050	0.027	0.050	0.20	1	1.00	U
Pyrene	< 0.050	0.025	0.050	0.20	1	1.00	U
Benzo (a) Anthracene	< 0.050	0.024	0.050	0.20	1	1.00	U
Chrysene	< 0.050	0.019	0.050	0.20	1	1.00	U
Benzo (k) Fluoranthene	< 0.050	0.023	0.050	0.20	1	1.00	U
Benzo (b) Fluoranthene	< 0.050	0.025	0.050	0.20	1	1.00	U
Benzo (a) Pyrene	< 0.050	0.036	0.050	0.20	1	1.00	U
Indeno (1,2,3-c,d) Pyrene	< 0.050	0.022	0.050	0.20	1	1.00	U
Dibenz (a,h) Anthracene	< 0.050	0.027	0.050	0.20	1	1.00	U
Benzo (g,h,i) Perylene	<0.050	0.022	0.050	0.20	1	1.00	U
Surrogate	Rec. (%)	Control Limi	its Qualifi	ers			
Nitrobenzene-d5	90	28-139					
2-Fluorobiphenyl	86	33-144					
p-Terphenyl-d14	87	23-160					



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

07/22/14 14-07-1466

Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch II
ES103	14-07-1466-1-A	07/21/14 08:55	Aqueous	GC/MS OO	07/22/14	07/22/14 22:46	140722L02
Comment(s): - Results were eval	uated to the MDL (DL), con	centrations >=	to the MDL (DL	_) but < RL (LOC	Q), if found, ar	e qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LO0</u>	<u>3</u>	<u>DF</u>	<u>Qualifiers</u>
Acetone	<10	6.0	10	20		1.00	U,ICH
Benzene	<0.50	0.14	0.50	1.0		1.00	U
Bromodichloromethane	< 0.50	0.21	0.50	5.0		1.00	U
Bromoform	<1.0	0.50	1.0	10		1.00	U
Bromomethane	<5.0	3.9	5.0	20		1.00	U
2-Butanone	<5.0	2.2	5.0	10		1.00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0		1.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0		1.00	U
Chloroethane	<5.0	2.3	5.0	10		1.00	U
Chloroform	<0.50	0.46	0.50	5.0		1.00	U
Chloromethane	<2.0	1.8	2.0	10		1.00	U,IJ
Dibromochloromethane	<0.50	0.25	0.50	1.0		1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10		1.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0		1.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0		1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0		1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0		1.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0		1.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0		1.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0		1.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0		1.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0		1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0		1.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0		1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0		1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10		1.00	U
Styrene	<0.50	0.17	0.50	1.0		1.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0		1.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0		1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0		1.00	U
Toluene	<0.50	0.24	0.50	1.0		1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0		1.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0		1.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0		1.00	U



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07/22/14 14-07-1466 **EPA 5030C** GC/MS / EPA 8260B

Units:

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Project: Red Hill LTM 112066

<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOQ	<u>DF</u>	<b>Qualifiers</b>
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U

Rec. (%)	Control Limits	<b>Qualifiers</b>
97	80-126	
93	80-134	
99	80-120	
99	88-112	
94	80-120	
	97 93 99 99	97 80-126 93 80-134 99 80-120 99 88-112





Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

07/22/14 14-07-1466

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES104	14-07-1466-2-A	07/21/14 10:00	Aqueous	GC/MS OO	07/22/14	07/22/14 18:15	140722L025
Comment(s): - Results were evaluated	d to the MDL (DL), con	centrations >= t	o the MDL (DL	) but < RL (LOC	), if found, are	qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	LOD	LOC	<u>]</u>	<u>DF</u>	<u>Qualifiers</u>
Acetone	<10	6.0	10	20	1	1.00	U,ICH
Benzene	<0.50	0.14	0.50	1.0	1	1.00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0	1	1.00	U
Bromoform	<1.0	0.50	1.0	10	1	1.00	U
Bromomethane	<5.0	3.9	5.0	20	1	1.00	U
2-Butanone	<5.0	2.2	5.0	10	1	1.00	U
Carbon Tetrachloride	< 0.50	0.23	0.50	1.0	1	.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0	1	1.00	U
Chloroethane	<5.0	2.3	5.0	10	1	1.00	U
Chloroform	< 0.50	0.46	0.50	5.0	1	.00	U
Chloromethane	<2.0	1.8	2.0	10	1	.00	U,IJ
Dibromochloromethane	< 0.50	0.25	0.50	1.0	1	.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1	.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1	1.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1	1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1	1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0	1	1.00	U
1,1-Dichloroethane	< 0.50	0.28	0.50	5.0	1	.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0	1	1.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0	1	1.00	U
c-1,2-Dichloroethene	< 0.50	0.48	0.50	1.0	1	.00	U
t-1,2-Dichloroethene	< 0.50	0.37	0.50	1.0	1	.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1	1.00	U
c-1,3-Dichloropropene	< 0.50	0.25	0.50	1.0	1	.00	U
t-1,3-Dichloropropene	< 0.50	0.25	0.50	1.0	1	.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0	1	1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0	1	.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1	.00	U
Styrene	<0.50	0.17	0.50	1.0	1	1.00	U
1,1,1,2-Tetrachloroethane	< 0.50	0.40	0.50	1.0	1	1.00	U
1,1,2,2-Tetrachloroethane	< 0.50	0.41	0.50	1.0	1	1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0	1	1.00	U
Toluene	<0.50	0.24	0.50	1.0	1	1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1	1.00	U
1,1,1-Trichloroethane	< 0.50	0.30	0.50	5.0	1	1.00	U
Hexachloro-1,3-Butadiene	< 0.50	0.32	0.50	1.0	1	1.00	U



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Date Received: Work Order: Preparation: Method:

07/22/14 14-07-1466 EPA 5030C

Units:

GC/MS / EPA 8260B ug/L

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Project: Red Hill LTM 112066

Toluene-d8

Toluene-d8-TPPH

1,4-Bromofluorobenzene

<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<b>Qualifiers</b>
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	0.36	0.23	0.50	1.0	1.00	J
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	48	26	30	50	1.00	J
Surrogate	Rec. (%)	Control Limits	<u>Qualifiers</u>			
Dibromofluoromethane	99	80-126				
1,2-Dichloroethane-d4	93	80-134				

80-120

88-112

80-120

100

99





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Date Received:

07/22/14 14-07-1466

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES105	14-07-1466-3-A	07/21/14 10:55	Aqueous	GC/MS OO	07/22/14	07/22/14 23:13	140722L025
Comment(s): - Results were evaluated to	o the MDL (DL), con	centrations >= t	o the MDL (DL	_) but < RL (LOC	(a), if found, are	qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	LOD	LOC	<u>D</u>	<u>)F</u>	<u>Qualifiers</u>
Acetone	<10	6.0	10	20	1	.00	U,ICH
Benzene	<0.50	0.14	0.50	1.0	1	.00	U
Bromodichloromethane	< 0.50	0.21	0.50	5.0	1	.00	U
Bromoform	<1.0	0.50	1.0	10	1	.00	U
Bromomethane	<5.0	3.9	5.0	20	1	.00	U
2-Butanone	<5.0	2.2	5.0	10	1	.00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0	1	.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0	1	.00	U
Chloroethane	<5.0	2.3	5.0	10	1	.00	U
Chloroform	<0.50	0.46	0.50	5.0	1	.00	U
Chloromethane	<2.0	1.8	2.0	10	1	.00	U,IJ
Dibromochloromethane	<0.50	0.25	0.50	1.0	1	.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1	.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1	.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1	.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1	.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0	1	.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1	.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0	1	.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0	1	.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0	1	.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0	1	.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1	.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1	.00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1	.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0	1	.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0	1	.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1	.00	U
Styrene	<0.50	0.17	0.50	1.0	1	.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1	.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0	1	.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0	1	.00	U
Toluene	<0.50	0.24	0.50	1.0	1	.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1	.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0		.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0		.00	U



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07/22/14 14-07-1466 EPA 5030C GC/MS / EPA 8260B

Units:

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Project: Red Hill LTM	112066

Toluene-d8

Toluene-d8-TPPH

1,4-Bromofluorobenzene

<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	< 0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	< 0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	0.33	0.23	0.50	1.0	1.00	J
Methyl-t-Butyl Ether (MTBE)	< 0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	49	26	30	50	1.00	J
<u>Surrogate</u>	Rec. (%)	Control Limits	<u>Qualifiers</u>			
Dibromofluoromethane	99	80-126				
1,2-Dichloroethane-d4	93	80-134				

80-120

88-112

80-120

99

98





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Date Received:

07/22/14

Work Order: Preparation:

14-07-1466 EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES TRIP	14-07-1466-4-A	07/21/14 08:00	Aqueous	GC/MS OO	07/22/14	07/22/14 17:48	140722L025
Comment(s): - Results were evaluated t	o the MDL (DL), con	centrations >= t	o the MDL (DI	_) but < RL (LOC	(a), if found, ar	e qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>Q</u>	<u>DF</u>	<u>Qualifiers</u>
Acetone	6.5	6.0	10	20		1.00	J,ICH
Benzene	<0.50	0.14	0.50	1.0		1.00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0		1.00	U
Bromoform	<1.0	0.50	1.0	10		1.00	U
Bromomethane	<5.0	3.9	5.0	20		1.00	U
2-Butanone	<5.0	2.2	5.0	10		1.00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0		1.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0		1.00	U
Chloroethane	<5.0	2.3	5.0	10		1.00	U
Chloroform	<0.50	0.46	0.50	5.0		1.00	U
Chloromethane	<2.0	1.8	2.0	10		1.00	U,IJ
Dibromochloromethane	<0.50	0.25	0.50	1.0		1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10		1.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0		1.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0		1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0		1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0		1.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0		1.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0		1.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0		1.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0		1.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0		1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0		1.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0		1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0		1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10		1.00	U
Styrene	<0.50	0.17	0.50	1.0		1.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0		1.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0		1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0		1.00	U
Toluene	<0.50	0.24	0.50	1.0		1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0		1.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0		1.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0		1.00	U



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Date Received: Work Order: Preparation: Method:

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Project: Red	Hill LTM	112066
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<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<b>Qualifiers</b>
1,1,2-Trichloroethane	< 0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	< 0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	< 0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	< 0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U
<u>Surrogate</u>	Rec. (%)	Control Limits	<u>Qualifiers</u>			
Dibromofluoromethane	99	80-126				
1.2 Diablaraathana d1	00	00.404				

Surrogate	Rec. (%)	Control Limits
Dibromofluoromethane	99	80-126
1,2-Dichloroethane-d4	92	80-134
Toluene-d8	99	80-120
Toluene-d8-TPPH	99	88-112
1.4-Bromofluorobenzene	95	80-120



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Lab Sample

Kailua, HI 96734-2500

Client Sample Number

Date Received:

07/22/14

Work Order: Preparation:

14-07-1466 EPA 5030C

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Date/Time

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Project: Red Hill LTM 112066

Date/Time QC Batch ID Matrix Instrument Date

Olient Gample Number	Number	Collected	WIGHTA	motrament	Prepared	Analyzed	QO Baterrib
Method Blank	099-13-057-57	N/A	Aqueous	GC/MS OO	07/22/14	07/22/14 17:12	140722L025
Comment(s): - Results were evaluated to	o the MDL (DL), cor	ncentrations >= t	o the MDL (DL	_) but < RL (LOC	Q), if found, a	re qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>3</u>	<u>DF</u>	<u>Qualifiers</u>
Acetone	<10	6.0	10	20		1.00	U
Benzene	< 0.50	0.14	0.50	1.0		1.00	U
Bromodichloromethane	< 0.50	0.21	0.50	5.0		1.00	U
Bromoform	<1.0	0.50	1.0	10		1.00	U
Bromomethane	<5.0	3.9	5.0	20		1.00	U
2-Butanone	<5.0	2.2	5.0	10		1.00	U
Carbon Tetrachloride	< 0.50	0.23	0.50	1.0		1.00	U
Chlorobenzene	< 0.50	0.17	0.50	5.0		1.00	U
Chloroethane	<5.0	2.3	5.0	10		1.00	U
Chloroform	< 0.50	0.46	0.50	5.0		1.00	U
Chloromethane	<2.0	1.8	2.0	10		1.00	U
Dibromochloromethane	< 0.50	0.25	0.50	1.0		1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10		1.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0		1.00	U
1,2-Dichlorobenzene	< 0.50	0.46	0.50	1.0		1.00	U
1,3-Dichlorobenzene	< 0.50	0.40	0.50	1.0		1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0		1.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0		1.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0		1.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0		1.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0		1.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0		1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0		1.00	U
c-1,3-Dichloropropene	< 0.50	0.25	0.50	1.0		1.00	U
t-1,3-Dichloropropene	< 0.50	0.25	0.50	1.0		1.00	U
Ethylbenzene	< 0.50	0.14	0.50	1.0		1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0		1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10		1.00	U
Styrene	<0.50	0.17	0.50	1.0		1.00	U
1,1,1,2-Tetrachloroethane	< 0.50	0.40	0.50	1.0		1.00	U
1,1,2,2-Tetrachloroethane	< 0.50	0.41	0.50	1.0		1.00	U
Tetrachloroethene	< 0.50	0.39	0.50	5.0		1.00	U
Toluene	< 0.50	0.24	0.50	1.0		1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0		1.00	U
1,1,1-Trichloroethane	< 0.50	0.30	0.50	5.0		1.00	U
Hexachloro-1,3-Butadiene	< 0.50	0.32	0.50	1.0		1.00	U



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Date Received: Work Order: Preparation: Method:

<u>LOD</u>

<u>LOQ</u>

<u>DF</u>

07/22/14 14-07-1466 EPA 5030C GC/MS / EPA 8260B

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Qualifiers U U U U U U U U

Project: Red Hill LTM 112066

<u>Parameter</u>

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1,1,2-Trichloroethane	< 0.50	0.38	0.50	1.0	1.00
Trichloroethene	<0.50	0.37	0.50	1.0	1.00
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00
p/m-Xylene	<1.0	0.30	1.0	10	1.00
o-Xylene	<0.50	0.23	0.50	1.0	1.00
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00
Gasoline Range Organics	<30	26	30	50	1.00
Surrogate	Rec. (%)	Control Limits	<u>Qualifiers</u>		
Surrogate Dibromofluoromethane	Rec. (%)	Control Limits 80-126	Qualifiers		
	<del></del>		Qualifiers		
Dibromofluoromethane	98	80-126	Qualifiers		
Dibromofluoromethane 1,2-Dichloroethane-d4	98 92	80-126 80-134	<u>Qualifiers</u>		
Dibromofluoromethane 1,2-Dichloroethane-d4 Toluene-d8	98 92 99	80-126 80-134 80-120	<u>Qualifiers</u>		

<u>DL</u>

Result



Environmental Science International, Inc.

Date Received:

Work Order:

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Kailua, HI 96734-2500

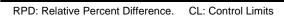
Preparation:

Method:

EPA 8015B (M)

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Quality Control Sample ID	Туре		Matrix	Ins	strument	Date Prepared	Date Ana	lyzed	MS/MSD Bat	ch Number
ES104	Sample		Aqueous	s GC	45	07/23/14	07/24/14	21:30	140723S11A	
ES104	Matrix Spike		Aqueous	s GC	45	07/23/14	07/24/14	20:17	140723S11A	
ES104	Matrix Spike	Duplicate	Aqueous	G GC	45	07/23/14	07/24/14	20:36	140723S11A	
Parameter	Sample Conc.	<u>Spike</u> Added	MS Conc.	<u>MS</u> %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
TPH as Diesel	1182	2000	3130	97	3204	101	55-133	2	0-30	





Environmental Science International, Inc.

Date Received:

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Kailua, HI 96734-2500

Preparation:

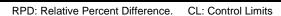
Method:

EPA 3005A Filt.

Method:

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Quality Control Sample ID	Туре		Matrix	In	nstrument	Date Prepared	Date Ana	lyzed	MS/MSD Bat	tch Number
ES104	Sample		Aqueou	s IC	CP/MS 04	07/23/14	07/24/14	19:48	140723S03	
ES104	Matrix Spike		Aqueou	s IC	CP/MS 04	07/23/14	07/24/14	19:39	140723S03	
ES104	Matrix Spike	Duplicate	Aqueou	s IC	CP/MS 04	07/23/14	07/24/14	19:41	140723S03	
Parameter	Sample Conc.	<u>Spike</u> Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Lead	ND	0.1000	0.1059	106	0.1128	113	80-120	6	0-20	





Environmental Science International, Inc. Da 354 Uluniu Street, Suite 304 Wo

Date Received: Work Order: Preparation: 07/22/14 14-07-1466 EPA 3510C

Method:

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EPA 8270C SIM PAHs

Project: Red Hill LTM 112066

Kailua, HI 96734-2500

Quality Control Sample ID	Туре		Matrix		Instrument	Date Prepare	ed Date Ana	lyzed	MS/MSD Ba	tch Number
ES104	Sample		Aqueou	ıs	GC/MS AAA	07/28/14	07/29/14	21:31	140728S01	
ES104	Matrix Spike		Aqueou	ıs	GC/MS AAA	07/28/14	07/29/14	13:52	140728S01	
ES104	Matrix Spike	Duplicate	Aqueou	ıs	GC/MS AAA	07/28/14	07/29/14	14:16	140728S01	
Parameter	Sample Conc.	<u>Spike</u> Added	MS Conc.	MS %Re	MSD ec. Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	<u>Qualifiers</u>
Naphthalene	70.83	2.000	59.97	0	64.80	0	21-133	8	0-25	3
2-Methylnaphthalene	19.64	2.000	21.30	83	18.58	0	21-140	14	0-25	3
1-Methylnaphthalene	25.19	2.000	25.26	4	24.73	0	20-140	2	0-25	3
Acenaphthylene	ND	2.000	1.370	68	1.495	75	33-145	9	0-25	
Acenaphthene	0.5154	2.000	1.776	63	1.936	71	49-121	9	0-25	
Fluorene	0.2363	2.000	1.536	65	1.689	73	59-121	10	0-25	
Phenanthrene	ND	2.000	1.313	66	1.517	76	54-120	14	0-25	
Anthracene	ND	2.000	1.421	71	1.551	78	27-133	9	0-25	
Fluoranthene	ND	2.000	1.282	64	1.431	72	26-137	11	0-25	
Pyrene	ND	2.000	1.192	60	1.344	67	18-168	12	0-25	
Benzo (a) Anthracene	ND	2.000	1.252	63	1.416	71	33-143	12	0-25	
Chrysene	ND	2.000	1.298	65	1.459	73	17-168	12	0-25	
Benzo (k) Fluoranthene	ND	2.000	1.137	57	1.284	64	24-159	12	0-25	
Benzo (b) Fluoranthene	ND	2.000	1.117	56	1.282	64	24-159	14	0-25	
Benzo (a) Pyrene	ND	2.000	1.181	59	1.348	67	17-163	13	0-25	
Indeno (1,2,3-c,d) Pyrene	ND	2.000	1.190	60	1.329	66	10-171	11	0-25	
Dibenz (a,h) Anthracene	ND	2.000	1.195	60	1.342	67	10-219	12	0-25	
Benzo (g,h,i) Perylene	ND	2.000	1.252	63	1.389	69	10-227	10	0-25	

GC/MS / EPA 8260B



#### **Quality Control - Spike/Spike Duplicate**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1466

Kailua, HI 96734-2500

Date Received:

Preparation:

EPA 5030C

Method:

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Quality Control Sample ID	Туре		Matrix	Ins	strument	Date Prepared	Date Ana	lyzed	MS/MSD Ba	tch Number
ES104	Sample		Aqueous	GC	C/MS OO	07/22/14	07/22/14	18:15	140722S011	
ES104	Matrix Spike		Aqueous	G	C/MS OO	07/22/14	07/22/14	18:42	140722S011	
ES104	Matrix Spike I	Duplicate	Aqueous	G	C/MS OO	07/22/14	07/22/14	19:10	140722S011	ı
Parameter	<u>Sample</u> <u>Conc.</u>	Spike Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Acetone	ND	50.00	68.38	137	67.89	136	40-140	1	0-20	
Benzene	ND	50.00	48.75	97	47.76	96	80-120	2	0-20	
Bromobenzene	ND	50.00	49.17	98	49.04	98	75-125	0	0-20	
Bromochloromethane	ND	50.00	51.02	102	50.51	101	65-135	1	0-20	
Bromodichloromethane	ND	50.00	50.05	100	48.80	98	75-120	3	0-20	
Bromoform	ND	50.00	49.04	98	47.81	96	70-130	3	0-20	
Bromomethane	ND	50.00	45.28	91	41.18	82	30-145	9	0-20	
2-Butanone	ND	50.00	51.87	104	50.83	102	30-150	2	0-20	
n-Butylbenzene	3.292	50.00	55.43	104	54.21	102	70-135	2	0-20	
sec-Butylbenzene	3.056	50.00	56.01	106	54.93	104	70-125	2	0-20	
tert-Butylbenzene	ND	50.00	53.19	106	52.66	105	70-130	1	0-20	
Carbon Disulfide	ND	50.00	37.91	76	37.81	76	35-160	0	0-20	
Carbon Tetrachloride	ND	50.00	46.80	94	46.16	92	65-140	1	0-20	
Chlorobenzene	ND	50.00	50.45	101	50.22	100	80-120	0	0-20	
Chloroethane	ND	50.00	42.16	84	42.32	85	60-135	0	0-20	
Chloroform	ND	50.00	47.60	95	47.54	95	65-135	0	0-20	
Chloromethane	ND	50.00	38.39	77	39.39	79	40-125	3	0-20	
2-Chlorotoluene	ND	50.00	51.15	102	50.66	101	75-125	1	0-20	
4-Chlorotoluene	ND	50.00	50.34	101	48.92	98	75-130	3	0-20	
Dibromochloromethane	ND	50.00	51.29	103	50.73	101	60-135	1	0-20	
1,2-Dibromo-3-Chloropropane	ND	50.00	44.35	89	44.28	89	50-130	0	0-20	
1,2-Dibromoethane	ND	50.00	48.68	97	48.61	97	80-120	0	0-20	
Dibromomethane	ND	50.00	50.83	102	50.24	100	75-125	1	0-20	
1,2-Dichlorobenzene	ND	50.00	50.78	102	50.20	100	70-120	1	0-20	
1,3-Dichlorobenzene	ND	50.00	51.20	102	50.63	101	75-125	1	0-20	
1,4-Dichlorobenzene	ND	50.00	49.14	98	48.08	96	75-125	2	0-20	
Dichlorodifluoromethane	ND	50.00	52.63	105	51.58	103	30-155	2	0-20	
1,1-Dichloroethane	ND	50.00	47.19	94	47.06	94	70-135	0	0-20	
1,2-Dichloroethane	ND	50.00	45.56	91	45.19	90	70-130	1	0-20	
1,1-Dichloroethene	ND	50.00	47.66	95	46.67	93	70-130	2	0-20	
c-1,2-Dichloroethene	ND	50.00	52.88	106	52.56	105	70-125	1	0-20	
t-1,2-Dichloroethene	ND	50.00	51.07	102	50.78	102	60-140	1	0-20	
1,2-Dichloropropane	ND	50.00	49.03	98	49.83	100	75-125	2	0-20	
1,3-Dichloropropane	ND	50.00	47.89	96	47.75	95	75-125	0	0-20	
2,2-Dichloropropane	ND	50.00	45.57	91	46.00	92	70-135	1	0-20	
,	<del>-</del>				. 5.00	<del>-</del> -		-		



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received:

14-07-1466

07/22/14

Work Order: Preparation:

EPA 5030C

Method:

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<u>Parameter</u>	Sample Conc.	<u>Spike</u> Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	<u>Qualifiers</u>
1,1-Dichloropropene	ND	50.00	47.94	96	48.39	97	75-130	1	0-20	
c-1,3-Dichloropropene	ND	50.00	52.46	105	52.70	105	70-130	0	0-20	
t-1,3-Dichloropropene	ND	50.00	51.27	103	51.49	103	55-140	0	0-20	
Ethylbenzene	ND	50.00	49.73	99	49.35	99	75-125	1	0-20	
2-Hexanone	ND	50.00	51.36	103	50.46	101	55-130	2	0-20	
Isopropylbenzene	ND	50.00	54.76	110	54.15	108	75-125	1	0-20	
p-Isopropyltoluene	ND	50.00	47.57	95	46.54	93	75-130	2	0-20	
Methylene Chloride	ND	50.00	50.87	102	50.51	101	55-140	1	0-20	
4-Methyl-2-Pentanone	ND	50.00	48.77	98	49.75	100	60-135	2	0-20	
Naphthalene	63.90	50.00	115.2	103	113.4	99	55-140	2	0-20	
n-Propylbenzene	4.683	50.00	56.47	104	55.41	101	70-130	2	0-20	
Styrene	ND	50.00	51.33	103	50.88	102	65-135	1	0-20	
1,1,1,2-Tetrachloroethane	ND	50.00	48.71	97	48.14	96	80-130	1	0-20	
1,1,2,2-Tetrachloroethane	ND	50.00	51.97	104	51.20	102	65-130	1	0-20	
Tetrachloroethene	ND	50.00	43.72	87	42.36	85	45-150	3	0-20	
Toluene	ND	50.00	48.84	98	47.83	96	75-120	2	0-20	
1,2,3-Trichlorobenzene	ND	50.00	50.43	101	50.15	100	55-140	1	0-20	
1,2,4-Trichlorobenzene	ND	50.00	50.65	101	49.25	99	65-135	3	0-20	
1,1,1-Trichloroethane	ND	50.00	46.97	94	46.67	93	65-130	1	0-20	
Hexachloro-1,3-Butadiene	ND	50.00	46.84	94	45.73	91	50-140	2	0-20	
1,1,2-Trichloro-1,2,2-Trifluoroethane	ND	50.00	45.86	92	44.56	89	80-130	3	0-20	
1,1,2-Trichloroethane	ND	50.00	50.20	100	49.27	99	75-125	2	0-20	
Trichloroethene	ND	50.00	48.36	97	46.73	93	70-125	3	0-20	
Trichlorofluoromethane	ND	50.00	48.95	98	47.58	95	60-145	3	0-20	
1,2,3-Trichloropropane	ND	50.00	46.90	94	47.05	94	75-125	0	0-20	
1,2,4-Trimethylbenzene	ND	50.00	51.25	103	50.42	101	75-130	2	0-20	
1,3,5-Trimethylbenzene	ND	50.00	51.97	104	51.17	102	75-130	2	0-20	
Vinyl Acetate	ND	50.00	58.38	117	57.09	114	80-120	2	0-20	
Vinyl Chloride	ND	50.00	42.41	85	41.82	84	50-145	1	0-20	
p/m-Xylene	ND	100.0	98.61	99	97.67	98	75-130	1	0-20	
o-Xylene	ND	50.00	52.60	105	51.94	104	80-120	1	0-20	
Methyl-t-Butyl Ether (MTBE)	ND	50.00	47.14	94	47.64	95	65-125	1	0-20	
Tert-Butyl Alcohol (TBA)	ND	250.0	236.1	94	230.4	92	46-154	2	0-35	
Diisopropyl Ether (DIPE)	ND	50.00	49.90	100	50.26	101	81-123	1	0-20	
Ethyl-t-Butyl Ether (ETBE)	ND	50.00	50.25	100	50.91	102	74-122	1	0-20	
Tert-Amyl-Methyl Ether (TAME)	ND	50.00	49.17	98	49.83	100	76-124	1	0-20	
Ethanol	ND	500.0	514.2	103	520.6	104	60-138	1	0-35	



#### **Quality Control - PDS**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1466

Kailua, HI 96734-2500

Preparation:

Method:

EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 1 of 1

Quality Control Sample ID	Туре	N	//atrix	Instrument	Date Prepared	Date Analyzed	PDS/PDSD Batch Number	
ES104	Sample	A	Aqueous	ICP/MS 04	07/23/14 00:00	07/24/14 19:48	140723S03	
ES104	PDS	A	Aqueous	ICP/MS 04	07/23/14 00:00	07/24/14 19:43	140723S03	
<u>Parameter</u>		Sample Conc.	Spike Added	PDS Conc	. PDS %Re	ec. %Rec. C	<u>Qualifiers</u>	
Lead		ND	0.1000	0.1023	102	75-125		





#### **Quality Control - LCS/LCSD**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1466

Kailua, HI 96734-2500

Preparation:

Method:

EPA 8015B (M)

Project: Red Hill LTM 112066 Page 1 of 5

Quality Control Sample ID	Type	Mat	trix	Instrument	Date Pre	pared Date	Analyzed	LCS/LCSD Ba	tch Number
099-15-516-158	LCS	Aqı	ueous	GC 45	07/23/14	07/24	4/14 18:30	140723B11A	
099-15-516-158	LCSD	Aqı	ueous	GC 45 07/23/		07/24	4/14 18:48	140723B11A	
Parameter	Spike Added	LCS Conc.	LCS %Rec.	LCSD Conc.	LCSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
TPH as Diesel	2000	2031	102	2018	101	60-132	1	0-11	





#### **Quality Control - LCS**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1466

Kailua, HI 96734-2500

Preparation:

Method:

EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 2 of 5

Quality Control Sample ID Type		Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number
099-14-497-87	LCS	Aqueous	ICP/MS 04	07/23/14	07/24/14 19:32	140723L03D
<u>Parameter</u>		Spike Added	Conc. Recovere	ed LCS %Re	ec. %Rec	. CL Qualifiers
Lead		0.1000	0.09175	92	80-12	0



#### **Quality Control - LCS**

Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Benzo (g,h,i) Perylene

Project: Red Hill LTM 112066

Date Received:

07/22/14 14-07-1466

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Page 3 of 5

Quality Control Sample ID	Туре	Matrix	Instrument	Date Prepared	Date Analyzed LCS B	atch Number
099-15-148-52	LCS	Aqueous	GC/MS AAA	07/28/14	07/29/14 13:28 14072	3L01
Parameter Parameter		Spike Added	Conc. Recove	ered LCS %Red	c. %Rec. CL	Qualifiers
Naphthalene		2.000	1.361	68	21-133	
2-Methylnaphthalene		2.000	1.230	61	21-140	
1-Methylnaphthalene		2.000	1.226	61	20-140	
Acenaphthylene		2.000	1.167	58	33-145	
Acenaphthene		2.000	1.271	64	55-121	
Fluorene		2.000	1.315	66	59-121	
Phenanthrene		2.000	1.379	69	54-120	
Anthracene		2.000	1.393	70	27-133	
Fluoranthene		2.000	1.385	69	26-137	
Pyrene		2.000	1.333	67	45-129	
Benzo (a) Anthracene		2.000	1.343	67	33-143	
Chrysene		2.000	1.447	72	17-168	
Benzo (k) Fluoranthene		2.000	1.265	63	24-159	
Benzo (b) Fluoranthene		2.000	1.294	65	24-159	
Benzo (a) Pyrene		2.000	1.273	64	17-163	
Indeno (1,2,3-c,d) Pyrene		2.000	1.408	70	25-175	
Dibenz (a,h) Anthracene		2.000	1.350	67	25-175	

1.500

75

25-157

2.000



#### **Quality Control - LCS/LCSD**

Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received: Work Order:

07/22/14 14-07-1466

Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Page 4 of 5

Quality Control Sample ID Type		Ma	trix	Instrument	Date Pre	pared Date	Analyzed	LCS/LCSD Ba	atch Number
099-13-057-57	LCS	Aqı	ueous	GC/MS OO	07/22/14	07/22	2/14 15:40	140722L025	
099-13-057-57	LCSD	Aqı	ueous	GC/MS OO	07/22/14	07/22	2/14 16:07	140722L025	
Parameter	Spike Added	LCS Conc.	LCS %Rec.	LCSD Conc.	LCSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Acetone	50.00	88.82	178	N/A	N/A	40-140	N/A	0-20	Χ
Benzene	50.00	49.88	100	N/A	N/A	80-120	N/A	0-20	
Bromobenzene	50.00	49.40	99	N/A	N/A	75-125	N/A	0-20	
Bromochloromethane	50.00	50.20	100	N/A	N/A	65-130	N/A	0-20	
Bromodichloromethane	50.00	49.69	99	N/A	N/A	75-120	N/A	0-20	
Bromoform	50.00	50.09	100	N/A	N/A	70-130	N/A	0-20	
Bromomethane	50.00	41.76	84	N/A	N/A	30-145	N/A	0-20	
2-Butanone	50.00	64.52	129	N/A	N/A	30-150	N/A	0-20	
n-Butylbenzene	50.00	54.15	108	N/A	N/A	70-135	N/A	0-20	
sec-Butylbenzene	50.00	55.17	110	N/A	N/A	70-125	N/A	0-20	
tert-Butylbenzene	50.00	54.15	108	N/A	N/A	70-130	N/A	0-20	
Carbon Disulfide	50.00	40.26	81	N/A	N/A	35-160	N/A	0-20	
Carbon Tetrachloride	50.00	49.66	99	N/A	N/A	65-140	N/A	0-20	
Chlorobenzene	50.00	51.79	104	N/A	N/A	80-120	N/A	0-20	
Chloroethane	50.00	42.88	86	N/A	N/A	60-135	N/A	0-20	
Chloroform	50.00	48.16	96	N/A	N/A	65-135	N/A	0-20	
Chloromethane	50.00	39.99	80	N/A	N/A	40-125	N/A	0-20	
2-Chlorotoluene	50.00	51.99	104	N/A	N/A	75-125	N/A	0-20	
4-Chlorotoluene	50.00	51.09	102	N/A	N/A	75-130	N/A	0-20	
Dibromochloromethane	50.00	50.95	102	N/A	N/A	60-135	N/A	0-20	
1,2-Dibromo-3-Chloropropane	50.00	47.49	95	N/A	N/A	50-130	N/A	0-20	
1,2-Dibromoethane	50.00	48.66	97	N/A	N/A	80-120	N/A	0-20	
Dibromomethane	50.00	50.64	101	N/A	N/A	75-125	N/A	0-20	
1,2-Dichlorobenzene	50.00	51.35	103	N/A	N/A	70-120	N/A	0-20	
1,3-Dichlorobenzene	50.00	52.64	105	N/A	N/A	75-125	N/A	0-20	
1,4-Dichlorobenzene	50.00	49.60	99	N/A	N/A	75-125	N/A	0-20	
Dichlorodifluoromethane	50.00	55.16	110	N/A	N/A	30-155	N/A	0-20	
1,1-Dichloroethane	50.00	48.68	97	N/A	N/A	70-135	N/A	0-20	
1,2-Dichloroethane	50.00	45.73	91	N/A	N/A	70-130	N/A	0-20	
1,1-Dichloroethene	50.00	50.18	100	N/A	N/A	70-130	N/A	0-20	
c-1,2-Dichloroethene	50.00	54.08	108	N/A	N/A	70-125	N/A	0-20	
t-1,2-Dichloroethene	50.00	52.36	105	N/A	N/A	60-140	N/A	0-20	
1,2-Dichloropropane	50.00	49.64	99	N/A	N/A	75-125	N/A	0-20	
1,3-Dichloropropane	50.00	47.76	96	N/A	N/A	75-125	N/A	0-20	
2,2-Dichloropropane	50.00	49.18	98	N/A	N/A	70-135	N/A	0-20	

50.00

50.36

1,1-Dichloropropene

N/A

N/A

75-130

N/A

0-20



#### **Quality Control - LCS/LCSD**

Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received: Work Order:

Work Order:
Preparation:
Method:

07/22/14 14-07-1466 EPA 5030C

GC/MS / EPA 8260B

Page 5 of 5

<u>Parameter</u>	Spike Added	LCS Conc.	LCS %Rec.	LCSD Conc.	LCSD %Rec.	%Rec. CL	<u>RPD</u>	RPD CL	<u>Qualifiers</u>
c-1,3-Dichloropropene	50.00	53.61	107	N/A	N/A	70-130	N/A	0-20	
t-1,3-Dichloropropene	50.00	53.53	107	N/A	N/A	55-140	N/A	0-20	
Ethylbenzene	50.00	51.23	102	N/A	N/A	75-125	N/A	0-20	
2-Hexanone	50.00	57.37	115	N/A	N/A	55-130	N/A	0-20	
Isopropylbenzene	50.00	53.77	108	N/A	N/A	75-125	N/A	0-20	
p-Isopropyltoluene	50.00	49.20	98	N/A	N/A	75-130	N/A	0-20	
Methylene Chloride	50.00	51.33	103	N/A	N/A	55-140	N/A	0-20	
4-Methyl-2-Pentanone	50.00	53.25	107	N/A	N/A	60-135	N/A	0-20	
Naphthalene	50.00	50.18	100	N/A	N/A	55-140	N/A	0-20	
n-Propylbenzene	50.00	53.12	106	N/A	N/A	70-130	N/A	0-20	
Styrene	50.00	52.08	104	N/A	N/A	65-135	N/A	0-20	
1,1,1,2-Tetrachloroethane	50.00	48.97	98	N/A	N/A	80-130	N/A	0-20	
1,1,2,2-Tetrachloroethane	50.00	50.50	101	N/A	N/A	65-130	N/A	0-20	
Tetrachloroethene	50.00	51.35	103	N/A	N/A	45-150	N/A	0-20	
Toluene	50.00	49.99	100	N/A	N/A	75-120	N/A	0-20	
1,2,3-Trichlorobenzene	50.00	50.65	101	N/A	N/A	55-140	N/A	0-20	
1,2,4-Trichlorobenzene	50.00	50.31	101	N/A	N/A	65-135	N/A	0-20	
1,1,1-Trichloroethane	50.00	48.83	98	N/A	N/A	65-130	N/A	0-20	
Hexachloro-1,3-Butadiene	50.00	48.70	97	N/A	N/A	50-140	N/A	0-20	
1,1,2-Trichloro-1,2,2-Trifluoroethane	50.00	48.22	96	N/A	N/A	80-130	N/A	0-20	
1,1,2-Trichloroethane	50.00	49.34	99	N/A	N/A	75-125	N/A	0-20	
Trichloroethene	50.00	51.40	103	N/A	N/A	70-125	N/A	0-20	
Trichlorofluoromethane	50.00	50.79	102	N/A	N/A	60-145	N/A	0-20	
1,2,3-Trichloropropane	50.00	48.87	98	N/A	N/A	75-125	N/A	0-20	
1,2,4-Trimethylbenzene	50.00	52.46	105	N/A	N/A	75-130	N/A	0-20	
1,3,5-Trimethylbenzene	50.00	52.49	105	N/A	N/A	75-130	N/A	0-20	
Vinyl Acetate	50.00	56.14	112	N/A	N/A	80-120	N/A	0-20	
Vinyl Chloride	50.00	43.53	87	N/A	N/A	50-145	N/A	0-20	
p/m-Xylene	100.0	101.3	101	N/A	N/A	75-130	N/A	0-20	
o-Xylene	50.00	53.13	106	N/A	N/A	80-120	N/A	0-20	
Methyl-t-Butyl Ether (MTBE)	50.00	47.94	96	N/A	N/A	65-125	N/A	0-20	
Tert-Butyl Alcohol (TBA)	250.0	229.7	92	N/A	N/A	46-154	N/A	0-25	
Diisopropyl Ether (DIPE)	50.00	49.67	99	N/A	N/A	81-123	N/A	0-20	
Ethyl-t-Butyl Ether (ETBE)	50.00	50.76	102	N/A	N/A	74-122	N/A	0-20	
Tert-Amyl-Methyl Ether (TAME)	50.00	49.96	100	N/A	N/A	76-124	N/A	0-20	
Ethanol	500.0	497.8	100	N/A	N/A	60-138	N/A	0-25	
Gasoline Range Organics	1000	1036	104	1019	102	80-120	2	0-20	



#### **Sample Analysis Summary Report**

Work Order: 14-07-1466				Page 1 of 1
Method	<u>Extraction</u>	Chemist ID	Instrument	Analytical Location
EPA 6020	EPA 3005A Filt.	598	ICP/MS 04	1
EPA 8015B (M)	EPA 3510C	628	GC 45	1
EPA 8270C SIM PAHs	EPA 3510C	923	GC/MS AAA	1
GC/MS / EPA 8260B	EPA 5030C	849	GC/MS OO	2

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841 Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841



#### **Glossary of Terms and Qualifiers**

Work Order: 14-07-1466 Page 1 of 1

Qualifiers	<u>Definition</u>
*	See applicable analysis comment.
<	Less than the indicated value.
>	Greater than the indicated value.
1	Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification.
2	Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification.
3	Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control.
4	The MS/MSD RPD was out of control due to suspected matrix interference.
5	The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference.
6	Surrogate recovery below the acceptance limit.
7	Surrogate recovery above the acceptance limit.
В	Analyte was present in the associated method blank.
BU	Sample analyzed after holding time expired.
BV	Sample received after holding time expired.
DL	The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence.
E	Concentration exceeds the calibration range.
ET	Sample was extracted past end of recommended max. holding time.
HD	The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
ICH	Initial calibration verification recovery is above the control limit for this analyte.
ICJ	Initial calibration verification recovery is below the control limit for this analyte.
IH	Calibration verification recovery is above the control limit for this analyte.
IJ	Calibration verification recovery is below the control limit for this analyte.
J	Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated.
JA	Analyte positively identified but quantitation is an estimate.
LOD	The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level.
LOQ	The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias.
Q	Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater.

- SG The sample extract was subjected to Silica Gel treatment prior to analysis.
- Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD).
- X % Recovery and/or RPD out-of-range.
- Z Analyte presence was not confirmed by second column or GC/MS analysis.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

# rn to Contents

#### Richard Villafania

(466)

From:

Ann Dang [ADang@esciencei.com] Monday, July 21, 2014 6:14 PM

Sent: To:

Richard Villafania

Cc:

Domonkos Feher; Traci Sylva; Jeff Hattemer

Subject:

112066 Red Hill shipment 7/21/14

Attachments:

COC 7-21-14.pdf

Hi Richard,

We shipped 2 coolers today, which you should receive tomorrow. The Fedex tracking is 804557917456 and 780042307911.

On the COC, the method VOCs (8260) should be marked. Corrections were made on the attached COC.

Thanks, Ann

Click here to report this email as spam.

CHAIN OF CUSTODY RECORD

99/1

Calscience Environmental Laboratories, Inc.

Return to Contents

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Calscience Environmental Laboratories, Inc. 

7440 Lincoln Way, Garden Grove, CA Other locations: Concord, San Luis Obi For courier service / sample contact sales@calscien

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CHAIN OF CUSTODY RECORD

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01/01/14 Revision

UNITED STATES US

TO SAMPLE CONTROL CALSCIENCE LABORATORIES 7440 LINCOLN WAY

**GARDEN GROVE CA 92841** 

FedEx Express

2 of 2 MPS# 7800 4230 7911 Mstr# 8045 5791 7456

TUE - 22 JUL AA STANDARD OVERNIGHT

0200

92841 CA-US SNA



ORIGIN ID:HNLA

SHIP DATE: c: UL14 ACTWGT: 49.3 LB CAD: /POS1501 DIMS: 24x13x14 IN

UNITED STATES US

BILL RECIPIENT

TO SAMPLE CONTROL **CALSCIENCE LABORATORIES** 7440 LINCOLN WAY

**GARDEN GROVE CA 92841** 

FedEx

1 of 2 THK# 0200 8045 5791 7456

VZ APVA

TUE - 22 JUL AA STANDARD OVERNIGHT

> 92841 CA-US SNA



Calscience

WORK ORDER #: 14-07- 1

# SAMPLE RECEIPT FORM

Cooler <u>i</u> of <u>2</u>

CLIENT: <u>Env. Science</u>	DATE:_	07/22/	<u>′ 14</u>
TEMPERATURE: Thermometer ID: SC1 (Criteria: 0.0 °C - 6.0 °C, not frozer Temperature	Blank ay of samp	☐ Sample	
CUSTODY SEALS INTACT:  □ Cooler □ □ No (Not Intact) □ Not Present □ Sample □ No (Not Intact) □ Not Present		Checked by Checked by	to the supplied and the state of the
SAMPLE CONDITION:	Yes	No	N/A
Chain-Of-Custody (COC) document(s) received with samples	.⊿′		
COC document(s) received complete	. <b>p</b> /		
<ul> <li>□ Collection date/time, matrix, and/or # of containers logged in based on sample labels.</li> <li>□ No analysis requested.</li> <li>□ Not relinquished.</li> <li>□ No date/time relinquished.</li> </ul>			
Sampler's name indicated on COC	Ø		
Sample container label(s) consistent with COC			
Sample container(s) intact and good condition			
Proper containers and sufficient volume for analyses requested			
Analyses received within holding time			
Aqueous samples received within 15-minute holding time			
□ pH □ Residual Chlorine □ Dissolved Sulfides □ Dissolved Oxygen			Ø
Proper preservation noted on COC or sample container			
☐ Unpreserved vials received for Volatiles analysis			
Volatile analysis container(s) free of headspace	.⊿		
Tedlar bag(s) free of condensation	. 🗖		d
	s <sup>®</sup> □ <u>T</u> erra	aCores® □_	
Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve () □EnCore Aqueous: ☑VOA ☑VOAh □VOAna₂ □125AGB □125AGBh □125AGBp	y / Z ☑1AGB	_ □1AGB <b>na</b> ₂ □	]1AGB <b>s</b>
□500AGB Ø500AGJ □500AGJs □250AGB □250CGB □250CGBs	□1PB	□1PBna □	500PB
□250PB 250PBnup□125PB □125PBznna □100PJ □100PJna2 □			
Air: Dedlar Canister Other: Dedlar Trip Blank Lot#: 1404227  Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Er	A Labeleo ivelope	I/Checked by: Reviewed by:	1/6



## Calscience

WORK ORDER #: 14-07- 4 6 6

# SAMPLE RECEIPT FORM Cooler 2 of 2

CLIENT: Env. Science	DATE: _	07/22/	/ 14
TEMPERATURE: Thermometer ID: SC1 (Criteria: 0.0 °C – 6.0 °C, not froze	n except se	ediment/tissue	e)
Temperature 2 • 4 °C - 0.3 °C (CF) = 2 • ) °C	Blank	☐ Sample	)
☐ Sample(s) outside temperature criteria (PM/APM contacted by:)		·	
☐ Sample(s) outside temperature criteria but received on ice/chilled on same d	ay of sampl	ina	
☐ Received at ambient temperature, placed on ice for transport by Co		19•	
	Juliel.	Checked by	15
Ambient Temperature: ☐ Air ☐ Filter		Checked by	y
CUSTODY SEALS INTACT:			
□ Cooler □ □ □ No (Not Intact) □ Not Present	□ N/A	Checked by	/: <u>15</u>
✓ Sample □ □ No (Not Intact) □ Not Present		Checked by	1
	i i i i i i i i i i i i i i i i i i i		
SAMPLE CONDITION:	Yes	No	N/A
Chain-Of-Custody (COC) document(s) received with samples	. 🖳		
COC document(s) received complete	. <b>Z</b>		
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.			
☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.			
Sampler's name indicated on COC	Ø		
Sample container label(s) consistent with COC	7		
Sample container(s) intact and good condition	<b>e</b> ,		
Proper containers and sufficient volume for analyses requested	4		
Analyses received within holding time	Ø		
Aqueous samples received within 15-minute holding time			
□ pH □ Residual Chlorine □ Dissolved Sulfides □ Dissolved Oxygen	. 🗆 ,		p'
Proper preservation noted on COC or sample container	. d		
☐ Unpreserved vials received for Volatiles analysis			
Volatile analysis container(s) free of headspace	. 🗆		Q
Tedlar bag(s) free of condensation	. 🗆		Ø
Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve() □EnCore	s <sup>®</sup> □Terra	Cores <sup>®</sup> □_	
Aqueous: □VOA □VOAh □VOAna₂ □125AGB □125AGBh □125AGBp			
□500AGB ☑500AGJ □500AGJs □250AGB □250CGB □250CGBs	□1PB	□1PB <b>na</b> □	500PB
□250PB			
Air: □Tedlar <sup>®</sup> □Canister Other: □ Trip Blank Lot#:			
Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Er Preservative: h: HCL n: HNO3 na2:Na2S2O3 na: NaOH p: H3PO4 s: H2SO4 u: Ultra-pure znna: ZnAc2+Na		Reviewed by: Scanned by:	dance.

#### **RAW DATA SHEET** FOR METHOD: EPA 8015B (M)

**WORK ORDER:** INSTRUMENT:

14-07-1466

GC 45

**EPA 3510C EXTRACTION:** 

D/T EXTRACTED: 2014-07-23 00:00

**ANALYZED BY:** 

628 D/T ANALYZED:

2014-07-24 21:13

**REVIEWED BY:** 

D/T REVIEWED:

**DATA FILE:** 

W:\GC 45\GC 45 DATA\2014\140724\14072415.D\14072415

1

**CLIENT SAMPLE NUMBER: ES103** 

LCS/MB BATCH: MS/MSD BATCH:

ug/L

140723B11A 140723S11A SAMPLE VOLUME / WEIGHT: FINAL VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

ADJUSTMENT RATIO TO PF:

人名英国国家 沙樓

0.50

COMMENT:

UNITS:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag.

INI. CONC DF

CONC

DL

LOD 12

LOQ 25

**QUAL** 

COMPOUND TPH as Diesel

13300

1.00

66.7

11

b

#### 

#### Area Percent Report

Data File Name :  $W:\GC_45\GC 45\DATA\2014\140724\14072415.D$ 

Vial Number : Vial 15

Page Number : 1
Operator : 682
Instrument : GC 45
Sample Name : 14-07-1466-1 Injection Number: 1 Sequence Line : 15

Instrument Method: C:\CHEM32\1\METHODS\ ->

3 VV 0.022

The property of the second

0.364

Acquired on : 24 Jul 14 9:13:38 PM

Report Created on: 25 Jul 14 06:11 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

Sig.	. 1	in W:\GC_45	5\GC 45 DATA\201	4\140724\ ->				
			Area					
		2 1/2	0.41	0	7.77	0 027	0.03	0

1 1	•		1	' '	•	
1	2.142	0.41	· 1	. <b>0 VV</b>	0.027	0.030
2	2.272	5.28		3 VV	0.026	0.392
3	2.299	1.55		1 VV	0.018	0.115
4	2.335	1.49		1 VV	0.020	0.110
5	2.382	1.89		1 VV	0.020	0.140
6	2.402	0.98		1 VV	0.015	0.073
7	2.435	1.27		1 VV	0.020	0.094
8	2.475	3.67	7	2 VV	0.025	0.273
9	2.513	5.66		4 VV	0.024	0.420
10	2.559	4.22		` 2 VV	0.026	0.314
11	2.593	2.54		2 VV	0.023	0.189
12	2.647	5.64		2 VV	0.039	0.419
13	2.688	5.74		2 VV	0.033	0.427
14	2.777	13.85		4 VV	0.047	1.030
15	2.808	10.68		5 VV	0.029	0.794
16	2.870	5.38		4 VV	0.020	0.400
17	2.886	3.77		4 VV	0.013	0.280
18	2.907	8.36		6 VV	0.020	0.622
19	2.938	11.55		8 VV	0.020	0.859
20	2.997	17.08		5 VV	0.044	1.269
21	3.063	10.17		4 VV	0.033	0.756
22	3.104	5.33		4 VV	0.020	0.396
23	3.123	2.22		3 VV	0.012	0.165
24	3.147	7.96		4 VV	0.028	0.592
25	3.220	20.24		11 VV	0.026	1.504
26	3.246	4.01		3 VV	0.021	0.298
27	3.286	5.37		3 VV	0.028	0.399
28	3.357	12.11		3 VV 🕒 🕾	0.046	0.900
29	3.387	7.69		·3 VV .	0.035	0.571
30	3.438	2.67		2 VV,	0.021	0.199
31	3.497	9.95	•	3 VV	0.044	0.739
32	3.549	9.61		3 VV 10.	0.047	0.714
33	3.616	8.98		3 VV	0.040	0.667
34	3.661	4.05		3 VV 🖘	0.022	0.301
35	3.699	4.70		2 VV	0.029	0.349
36	3.744	5.38		2 VV	0.033	0.400
37	3.772	3.19		2 VV ;,	0.023	0.237
38	3.806	4.31		2 VV	0.029	0.320
39	3.833	2.79		2 VV	0.027	0.208
40	3.872	2.04		2 VV	0.019	0.151
41	3.896	2.75		2 VV	0.021	0.204
42	3.912	2.87		2 VV	0.030	0.213
43	3.960	2.17		1 VV	0.021	0.161
44	3.986	5.24		2 VV	0.042	0.389
		4 00		2 777	0 000	0 264

4.90

45

4.058

#### Area Percent Report

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Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072415.D

Page Number : 2

Operator : 682 Vial Number : Vial 15

Instrument : GC 45 Injection Number : 1 Sample Name : 14-07-1466-1 Sequence Line : 15

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 24 Jul 14 9:13:38 PM

Report Created on: 25 Jul 14 06:11 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

Pk I	Ret Time	Area	Height	Peak	Width	Response %
				-		
46	4.099	2.43		1 VV	0.041	0.180
47	4.166	1.21		1 VV	0.030	0.090
48	4.218	1.17		1 VV	0.028	0.087
49	4.267	1.72		1 VV	0.032	0.128
50	4.307	0.61		$\nabla V^{'}$ 0	0.028	0.045
51	4.351	0.33		0 VV 0	0.025	0.024
52	4.390	0.41	٤ .	0 VV	0.024	0.030
53	4.424	1.06		1 VB	0.030	0.079
54	4.566	0.00	• •	0 BV	0.000	0.000
55	4.602	0.29		0 VB	. 0.021	0.021
56	4.681	0.85		1 BV	0.018	0.063
57	4.713	2.88		2 VV ,	0.024	0.214
58	4.748	0.54		0 VB	0.021	0.040
59	4.860	0.30		0 BV	0.023	0.022
60	4.910	0.61		0 VV	(5) 0.022	0.045
61	4.967	0.35	. ez	0 VB	0.020	0.026
62	5.028	0,26		0 BV	0.017	0.019
63	5.059	1.69		1 VV	0.022	0.126
64	5.124	0.06		0 VV	0.031	0.004
65	5.205	1.43		1 VV	0.023	0.106
66	5.259	0.58		0 VV	0.035	0.043
67	5.353	1065.17	81	1 VV	0.020	79.157

Total area = 1345.64

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#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072415.D

Page Number : 3

Sample Name

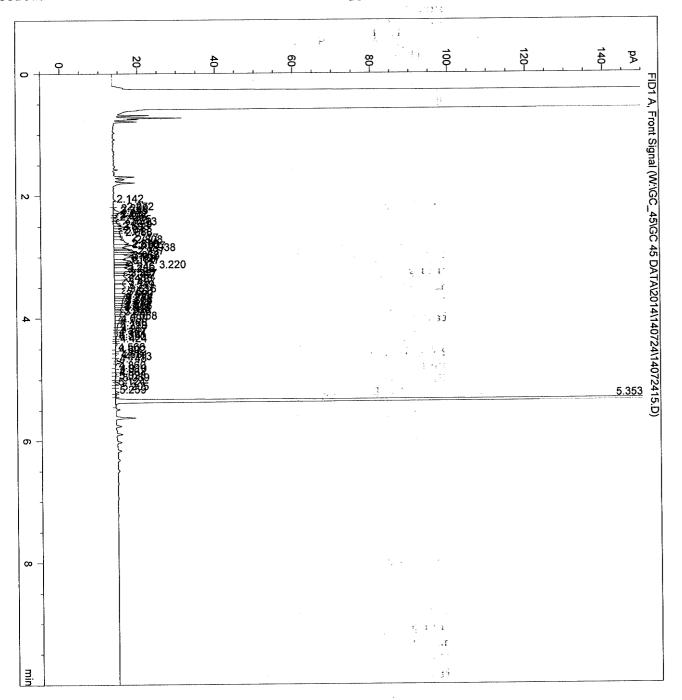
Operator : 682 Instrument : GC 4 Vial Number : Vial 15

Instrument Method: C:\CHEM32\1\METHODS\

Acquired on : 24 Jul 14 9:13:38 PM

Report Created on: 25 Jul 14 06:11 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies



# **RAW DATA SHEET** FOR METHOD: EPA 8015B (M)

WORK ORDER: **INSTRUMENT:** 

14-07-1466

GC 45

**EXTRACTION: EPA 3510C** 

2014-07-23 00:00 D/T EXTRACTED:

ANALYZED BY:

628

D/T ANALYZED: 2014-07-24 21:30

**REVIEWED BY:** D/T REVIEWED:

DATA FILE:

W:\GC 45\GC 45 DATA\2014\140724\14072416.D\14072416

**CLIENT SAMPLE NUMBER: ES104** 

LCS/MB BATCH: MS/MSD BATCH:

140723B11A 140723S11A ug/L

SAMPLE VOLUME / WEIGHT:

FINAL VOLUME / WEIGHT:

ADJUSTMENT RATIO TO PF:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

0.50

COMMENT:

UNITS:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag.

TPH as Diesel

INI. CONC

DF

CONC :

DL

LOD 12

LOQ 25

QUAL

COMPOUND

236000

1.00

1180

11

b

## 

#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072416.D

Vial Number : Vial 16

Page Number : 1
Operator : 682
Instrument : GC 45
Sample Name : 14-07-1466-2 Injection Number: 1 Sequence Line : 16

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 24 Jul 14 9:30:56 PM

Report Created on: 29 Jul 14 01:05 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

Sig. 1	in W:\GC_45\GC	45 DATA\203	L4\140724\	->			
Pk	Ret Time	Area	Height		Peak	Width	Response %
	-						
1	1.944	8.11		4	VV	0.030	0.138
2	1.994	26.82		16	VV	0.024	0.455
3	2.077	58.15		56	VV	0.016	0.987
4	2.107	28.67		27	VV .	0.016	0.487
5	2.136	35.34		22	VV	0.022	0.600
6	2.220	151.81		84	VV	0.028	2.578
7	2.271	52.19		29	VV	0.025	0.886
8	2.304	69.93		54	VV	0.020	1.187
9	2.333	21.11		18	VV.	0.017	0.358
10	2.371	69.23		42	ΔΔ.,	0.024	1.175
11	2.401	91.73		74	VV	0.018	1.558
12	2.436	62.19		42	VV	0.021	1.056
13	2.507	466.71		324	VV	0.021	7.925
14	2.559	38.11		28	VV .	0.019	0.647
15	2.596	79.47	á –	36	$\Delta \Delta_{p, q = f}$	0.028	1.349
16	2.616	36.59		31	VV	0.017	0.621
17	2.650	77.13	in the state of	40	. ΔΔ.,	0.029	1.310
18	2.690	96.26		44	VV	0.029	1.635
19	2.731	53.13		40	VV	0.020	0.902
20	2.781	281.08		146	VV	0.026	4.773
21	2.822	259.67		157		0.023	4.409
22	2.871	52.51			VV	0.016	0.892
23	2.905	189.91	•	83	VV	0.031	3.225
24	2.938	126.05		61	VV	0.028	2.140
25	2.998	179.68		67	VV	0.039	3.051
26		58.66		50	VV	0.017	0.996
27		195.20		104	.vv	0.026	3.314
28		132.36		68	$\nabla V$ .	0.029	2.248
29		103.41		47	VV	0.030	1.756
30		100.22		49	·VV	0.028	1.702
31		56.86		44	ΛΛ	0.018	0.965
32		75.74		42	VV	0.027	1.286
33		71.53		39	VV	0.026	1.215
34		132.16		51	<b>VV</b>	.% 0.036	2.244
35		56.69	•	51	νν :-	9 0.016	0.963
36		94.90	2	53	VV	0.025	1.611
37		47.97	3.4		<b>VV</b> ;5 &	cá 0.019	0.815
38		55.29			VV	0.021	0.939
39		106.30	4	43		. 0.032	1.805
40		80.63		43	VV	0.025	1.369
41		63.43	÷* ,		< <b>VV</b>	0.022	1.077
42		81.73		• • •	VV	0.029	1.388
43		117.41			VV	0.038	1.994
44		94.65	:		vv	0.038	1.607
45		38.76			. VV	0.021	0.658
1	2						

#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072416.D

Page Number : 2

Operator : 682 Vial Number : Vial 16

Instrument : GC 45 : Injection Number : 1
Sample Name : 14-07-1466-2 : Sequence Line : 16

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 24 Jul 14 9:30:56 PM

Report Created on: 29 Jul 14 01:05 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

CHAL	C ROVEDECII.			<b>.</b>	•	:	-
Pk	Ret Time	Area	Height		Peak	Width	Response %
							1 530
46	3.745	90.24		35	VV	0.034	1.532
47	3.771	64.83		32		0.029	1.101
48	3.827	76.08			VV	0.045	1.292
4.9	3.868	34.64			VV	0.024	0.588
50	3.907	38.12			VV	0.027	0.647
51	3.929	24.21			VV	0.018	0.411
52	3.953	25.33			VV	0.024	0.430
53	3.991	60.08			ΔΛ	0.043	1.020
54	4.056	34.15			VV	0.033	0.580
55	4.098	22.78		9	ŢVV	0.032	0.387
56	4.136	10.52	: •	8	VV	0.020	0.179
57	4.164	12.26	•	7	VV	0.025	0.208
58	4.196	13.83		,6	VV.	0.032	0.235
59	4.232	6.45		5	, VV	0.017	0.110
60	4.251	12.16		5	VV	0.038	0.206
61	4.307	11.66		4	$\Delta \Delta \Delta$	0.044	0.198
62	4.352	3.90		3	VV ,	0.018	0.066
63	4.384	6.84		3	VV	0.030	0.116
64	4.419	17.93	• :	3	VV	0.076	0.304
65	4.589	4.78		1.	, AA	0.053	0.081
66	4.706	11.05		4	ΔΛ.	0.036	0.188
67	4.800	1.13	erā .	1	$vv^a$	0.028	0.019
68	4.856	1.60		,1	VV	0.035	0.027
69	4.906	1.18		. 0	ΛΛ	0.031	0.020
70	4.961	1.18	<i>3</i>	1	VV	0.030	0.020
71	5.021	0.43		0	VV	0.021	0.007
72	5.052	1.69		1	VV	0.023	0.029
73	5.094	0.05		0	VV	0.016	0.001
74	5.197	1.35		1	BV	0.023	0.023
75	5.249	0.42		0	VV	0.026	0.007
76	5.344	922.91		653	<b>VV</b> :	0.023	15.671

Total area = 5889.27

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072416.D

Page Number : 3

Sample Name

: 682 Operator Instrument

: GC 45 : 14-07-1466-2

: Vial 16 Vial Number

Injection Number: 1 Sequence Line : 16

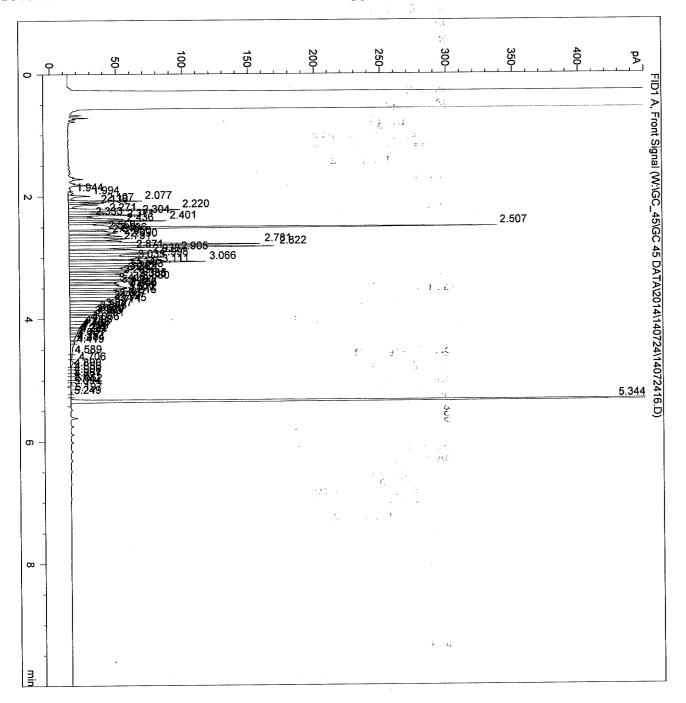
Instrument Method: C:\CHEM32\1\METHODS\

Acquired on : 24 Jul 14

9:30:56 PM 01:05 pm Report Created on: 29 Jul 14

Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies



# **RAW DATA SHEET** FOR METHOD: EPA 8015B (M)

33

**WORK ORDER: INSTRUMENT:** 

14-07-1466

GC 45

**EPA 3510C** 

**EXTRACTION:** D/T EXTRACTED:

2014-07-23 00:00

**ANALYZED BY:** 

D/T ANALYZED:

2014-07-24 21:48

628

**REVIEWED BY:** D/T REVIEWED:

**DATA FILE:** 

W:\GC 45\GC 45 DATA\2014\140724\14072417.D\14072417

E. 122 G

**CLIENT SAMPLE NUMBER: ES105** 

LCS/MB BATCH: MS/MSD BATCH:

140723B11A 140723S11A ug/L

SAMPLE VOLUME / WEIGHT: FINAL VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

ADJUSTMENT RATIO TO PF:

0.50

**COMMENT:** 

UNITS:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

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qualified with a "J" flag.

INI. CONC DF

-

CONC 1320

DL 11

LOD 12

QUAL LOQ

b

COMPOUND TPH as Diesel

265000. 1.00

31. 14 h

25

Page 4 of 4

# 

#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072417.D

Page Number : 1

Vial Number : Vial 17 : 682 Operator

Injection Number: 1 Instrument : GC 45 Sample Name : 14-07-1466-3 Sequence Line : 17

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 24 Jul 14 9:48:52 PM

Analysis Method : 8015B.MTH Report Created on: 29 Jul 14 01:05 pm

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

ig. 1 :	in W:\GC_45\G	C 45 DATA\2014	\140724\ ->	* A		
Pk	Ret Time	Area	Height	Peak	Width	Response %
						,
1	1.945	9.80		VV	0.030	0.146
2	1.994	31.86		VV	0.025	0.476
3	2.077	66.79		VV.	0.016	0.998
4	2.107	33.21		VV	0.016	0.496
5	2.136	41.20		VV	0.024	0.616
6	2.219	176.26		VV	0.029	2.633
7	2.271	60.30	the second secon	VV	0.030	0.901
8	2.304	82.33	62	VV	0.020	1.230
9	2.334	23.70		, VV	0.016	0.354
10	2.371	80.18	and the second s	νν΄	0.024	1.198
11	2.401	107.56		VV	0.018	1.607
12	2.437	71.67		VV	0.023	1.071
13	2.507	538.09	371	VV	0.021	8.038
14	2.560	47.85		νν , , , , , , , , , , , , , , , , , ,	0.021	0.715
15	2.597	87.08		VV	0.028	1.301
16	2.617	41.00	· ·	VV	0.017	0.612
17	2.653	85.64		VV	0.029	1.279
18	2.691	107.63	48	VV	0.031	1.608
19	2.731	58.32	44	VV	0.019	0.871
20	2.782	323.90	177	VV :	0.025	4.839
21	2.822	298.78	181	VV	0.023	4.463
22	2.906	264.71	89	VV	0.039	3.954
23	2.939	137.94	67	VV.	0.027	2.061
24	2.998	195.07	; 74	· V,V	0.038	2.914
25	3.037	68.68		VV .	0.017	1.026
26	3.067	216.73	121	VV	0.024	3.238
27	3.111	145.33		$\Delta \Delta$		2.171
28	3.146	113.40		<b>VV</b>		1.694
29	3.194	112.52	53	VV	0.028	1.681
30	3.223	79.27	46		0.022	1.184
31	3.250	58.65	45		0.022	0.876
32	3.283	69.30	4.3		0.023	1.035
33	3.333	135.78	56		0.032	2.028
34	3.355	75.60		VV	0.020	1.129
35	3.380	111.65		VV -	0.028	1.668
36	3.416	48.71		. VV :		0.728
37	3.441	66.84	40	) VV		0.998
38	3.487	100.77		VV	0.030	1.505
39	3.506	91.91		vv v	0.027	
40	3.551	68.65		S VV	0.021	1.026
41	3.575	89.78		VV	0.028	1.341
42	3.617	107.77		S VV	0.031	
43	3.644	34.70		γV	0.015	0.518
44	3.667	69.31		VV (	0.026	
45	3.689	59.20	35	VV.	0.025	0.884

#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072417.D

Page Number : 2

Vial Number : Vial 17 : 682 Operator

Injection Number: 1 : GC 45 Instrument Sequence Line : 17 Sample Name : 14-07-1466-3

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 24 Jul 14 9:48:52 PM

Analysis Method : 8015B.MTH Report Created on: 29 Jul 14 01:05 pm

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

Pk	Ret Time	Area	Height	Peak	Width	Response %
		08.51		•	0.015	0.411
46	3.721	27.51		31 VV	0.013	1.047
47	3.747	70.09		10 VV	0.023	
48	3.771	73.69		34 VV		
49	3.832	91.10		27 VV	0.043	
50	3.872	36.78		22 VV	0.027	
51	3.909	60.17		22 VV	0.036	
52	3.953	33.43	;	L9 VV	0.027	
53	3.992	65.23		20 VV	0.041	
54	4.058	37.63		15 VV	0.033	
55	4.100	35.70		LO VV	0.044	
56	4.167	15.18		8 VV	0.026	
57	4.197	14.10		7 VV	0.027	
58	4.229	9.65		6 VV	0.021	
59	4.254	12.68	2	6 VV	0.036	
60	4.303	11.50		4 VV	0.038	
61	4.349	6.21		4 VV	0.024	
62	4.386	7.68	<b>,</b> .	4 VV	0.028	
63	4.419	7.29		4 VV	0.029	
64	4.451	2.33	*	3, VV	0.015	
65	4.472	7.09		3 VV	0.035	
66	4.520	7.25		2 VV	0.048	
67	4.594	5.48		1 VV	0.061	
68	4.710	12.26		3 VV	0.052	
69	4.794	2.35		1 VV	0.040	
70	4.851	2.83		1 VV	0.042	
71	4.905	2.36		1 VV	0.042	
72	4.962	2.53		1 VV	0.044	
73	5.055	4.44		2 VV	0.036	
74	5.124	0.50		$\nabla V = 0$	0.018	
75	5.199	2.89		1 VV	0.034	
76	5.251	1.43	-	1 · VV	0.039	0.021
77	5.348	1129.37	. 8	60 VV	0.020	16.871
				. 3	1.0	

Total area = 6694.13

#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072417.D

Page Number : 3

Operator : 682 Vial Number : Vial 17

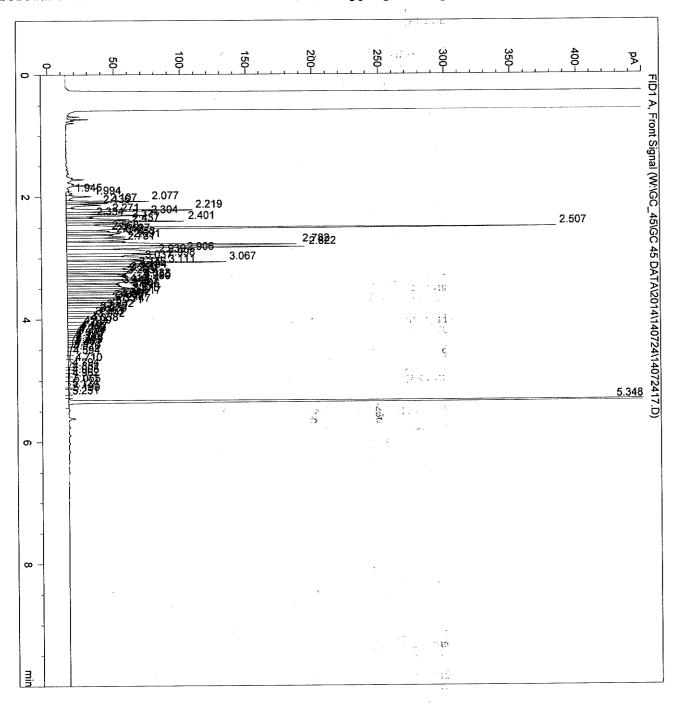
Instrument : GC 45 Injection Number : 1
Sample Name : 14-07-1466-3 Sequence Line : 17

Instrument Method: C:\CHEM32\1\METHODS\ -

Acquired on : 24 Jul 14 9:48:52 PM

Report Created on: 29 Jul 14 01:05 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies



#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072405.D

Page Number

: Vial 5 Vial Number

Operator Instrument : 682 : GC 45

Injection Number: 1

Sequence Line : 5

Sample Name

: MB 14072311/12

Instrument Method: C:\CHEM32\1\METHODS\

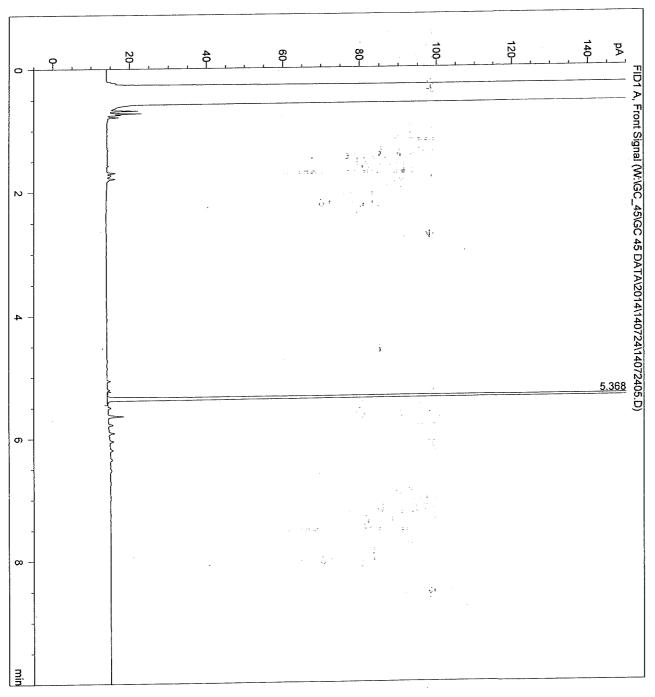
Acquired on

: 24 Jul 14

6:11:35 PM

Report Created on: 25 Jul 14 06:10 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies



#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072403.D

?age Number perator

: Vial 3 Vial Number

: 682 : GC 45

Injection Number: 1 Sequence Line

[nstrument Jample Name

: D400 C28 50 L041814D

Instrument Method: C:\CHEM32\1\METHODS\

Acquired on Report Created on: 25 Jul 14

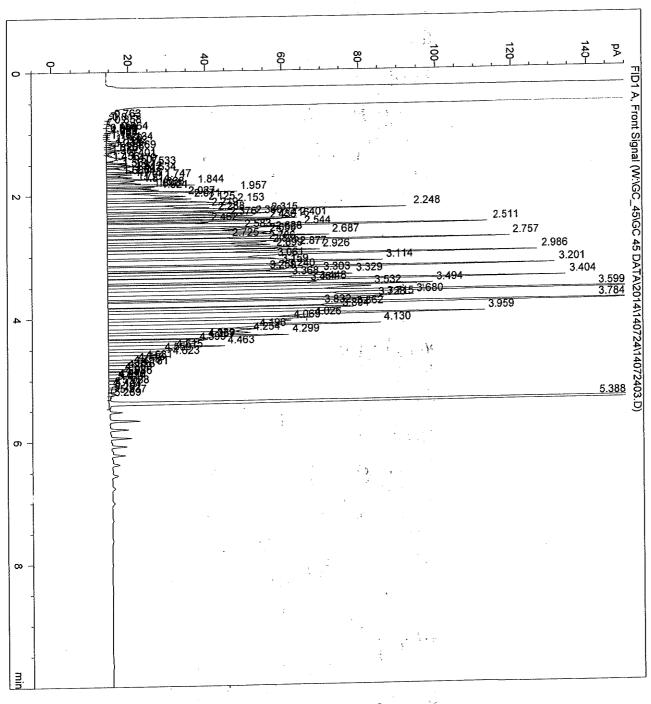
: 24 Jul 14

5:36:31 PM

Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

06:10 pm





#### Supplemental Report 1

# Calscience



# **WORK ORDER NUMBER: 14-07-1550**

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

**Analytical Report For** 

Client: Environmental Science International, Inc.

Client Project Name: Red Hill 112066

**Attention:** Robert Chong

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Richard Vellas

Approved for release on 07/30/2014 by: Richard Villafania

Project Manager



ResultLink >
Email your PM >

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.



# **Contents**

Client Project Name: Red Hill 112066 Work Order Number: 14-07-1550

WOIK OI	del Number. 14 07-1330	
1	Work Order Narrative	3
2	Client Sample Data	4
3	Quality Control Sample Data.  3.1 MS/MSD.  3.2 PDS/PDSD.  3.3 LCS/LCSD.	5 5 6 7
4	Sample Analysis Summary	8
5	Glossary of Terms and Qualifiers	9
6	Chain-of-Custody/Sample Receipt Form	10



#### **Work Order Narrative**

Work Order: 14-07-1550 Page 1 of 1

#### **Condition Upon Receipt:**

Samples were received under Chain-of-Custody (COC) on 07/23/14. They were assigned to Work Order 14-07-1550.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

#### **Holding Times:**

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

#### **Quality Control:**

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

#### **Additional Comments:**

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New\_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

#### **Subcontractor Information:**

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

U



Lead

### **Analytical Report**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

Units:

Units:

Units:

07/23/14

Work Order:

14-07-1550

Preparation:

N/A

Method:

Units:

Ug/L

Project: Red Hill 112066 Page 1 of 1

<0.0898

Client Sample I	Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES107UF		14-07-1550-5-A	07/22/14 09:45	Aqueous	ICP/MS 04	07/23/14	07/24/14 20:02	140723L02D
Comment(s):	- Results were evaluated	to the MDL (DL), con-	centrations >= t	o the MDL (D	L) but < RL (LC	Q), if found, are	qualified with a	"J" flag.
<u>Parameter</u>		<u>Resu</u>	<u>ılt</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Q</u>	<u>ualifiers</u>

Method Blank	099-16-094-424	N/A	Aqueous	ICP/MS 03	07/23/14	07/23/14 15:02	140723L02D
Comment(s):	- Results were evaluated to the MDL (DL), cor	ncentrations >=	to the MDL (DL	.) but < RL (LO	Q), if found, are	qualified with a	"J" flag.
<u>Parameter</u>	Res	<u>ult</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Q</u>	<u>tualifiers</u>
Lead	<0.0	898	1.00	0.0898	1.00	U	1

1.00

0.0898

1.00



RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



# **Quality Control - Spike/Spike Duplicate**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

Date Received:

07/23/14

Work Order:

14-07-1550

Preparation:

N/A

EPA 200.8

Project: Red Hill 112066 Page 1 of 1

Quality Control Sample ID	Туре		Matrix	Ins	strument	Date Prepared	Date Ana	lyzed	MS/MSD Bat	tch Number
14-07-1564-2	Sample		Aqueous	s IC	P/MS 03	07/23/14	07/23/14	15:15	140723S02	
14-07-1564-2	Matrix Spike		Aqueous	s IC	P/MS 03	07/23/14	07/23/14	15:09	140723S02	
14-07-1564-2	Matrix Spike I	Duplicate	Aqueous	s IC	P/MS 03	07/23/14	07/23/14	15:12	140723S02	
Parameter	Sample Conc.	<u>Spike</u> Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Lead	112.2	100.0	229.4	117	233.0	121	80-120	2	0-20	3



# **Quality Control - PDS**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

Date Received:

07/23/14

Work Order:

14-07-1550

Preparation:

N/A

EPA 200.8

Project: Red Hill 112066 Page 1 of 1

Quality Control Sample ID	Туре	N	Matrix	Instrument	Date Prepared	Date Analyzed	PDS/PDSD Batch Number
14-07-1564-2	Sample		Aqueous	ICP/MS 03	07/23/14 00:00	07/23/14 15:15	140723S02
14-07-1564-2	PDS		Aqueous	ICP/MS 03	07/23/14 00:00	07/29/14 14:31	140723S02
<u>Parameter</u>		Sample Conc.	Spike Added	PDS Conc.	PDS %Re	ec. %Rec. C	<u>CL</u> <u>Qualifiers</u>
Lead		112.2	500.0	640.0	106	75-125	





# **Quality Control - LCS**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

Date Received:

07/23/14

Work Order:

14-07-1550

Preparation:

N/A

EPA 200.8

Project: Red Hill 112066 Page 1 of 1

Quality Control Sample ID	Туре	Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number
099-16-094-424	LCS	Aqueous	ICP/MS 03	07/23/14	07/23/14 15:05	140723L02D
Parameter		Spike Added	Conc. Recovere	ed LCS %Re	ec. %Rec	. CL Qualifiers
Lead		100.0	95.43	95	80-120	0





# **Sample Analysis Summary Report**

Work Order: 14-07-1550				Page 1 of 1
Method	Extraction	Chemist ID	Instrument	Analytical Location
EPA 200.8	N/A	598	ICP/MS 04	1



#### **Glossary of Terms and Qualifiers**

Work Order: 14-07-1550 Page 1 of 1

Qualifiers	<u>Definition</u>
*	See applicable analysis comment.
<	Less than the indicated value.
>	Greater than the indicated value.
1	Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification.
2	Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification.
3	Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control.
4	The MS/MSD RPD was out of control due to suspected matrix interference.
5	The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference.
6	Surrogate recovery below the acceptance limit.
7	Surrogate recovery above the acceptance limit.
В	Analyte was present in the associated method blank.
BU	Sample analyzed after holding time expired.
BV	Sample received after holding time expired.
DL	The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence.
E	Concentration exceeds the calibration range.
ET	Sample was extracted past end of recommended max. holding time.
HD	The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
ICH	Initial calibration verification recovery is above the control limit for this analyte.
ICJ	Initial calibration verification recovery is below the control limit for this analyte.
IH	Calibration verification recovery is above the control limit for this analyte.
IJ	Calibration verification recovery is below the control limit for this analyte.
J	Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated.
JA	Analyte positively identified but quantitation is an estimate.
LOD	The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level.
LOQ	The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias.
Q	Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater.

- SG The sample extract was subjected to Silica Gel treatment prior to analysis.
- U Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD).
- X % Recovery and/or RPD out-of-range.
- Z Analyte presence was not confirmed by second column or GC/MS analysis.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

Calscience Environmental Laboratories, Inc. 7440 Lincoln Way, Garden Grove, CA 92841-1427。 (714) 895-5494 

Other locations: Concord, San Luis Obispo, Houston, and Corpus Christi For courier service / sample drop off information, contact sales@calscience.com or call us. LABORATORY CLIENT: ETN' IN INMENTAL Schence International

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CHAIN OF CUSTODY RECORD Page, Date

17.00°C CLIENT PROJECT NAME / NUMBER: RA KE

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PROJECT CONTACT

945 OLYST

STATE

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ADDRESS: 354 WILLING ST.

REQUESTED ANALYSES

SAMPLER(S): (PRINT) いかっているで

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Please check box or fill in blank as needed.

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STANDARD

5 DAYS

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☐ 48 HR

24 HR

GLOBAL ID

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COELT EDF SAME DAY

TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD");

TEL: 808-261-0740

CITY Kailud

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MIS 07S8 € 07S8 ☐ 2HA9

**LCB**2 (8085)

Pesticides (8081)

Oxygenates (8260)

BTEX / MTBE 🗆 8260 🗆

AOCs (85e0)

Lead

Page 10 of 15

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Time:

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Time:

Time:

01/01/14 Revision

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## Richard Villafania

From: Sent:

Ann Dang [ADang@esciencei.com] Tuesday, July 22, 2014 6:34 PM Richard Villafania

To:

Cc: Subject: Domonkos Feher; Traci Sylva; Jeff Hattemer

112066 Red Hill shipment 7-22-14

Attachments:

COC 7-22-14.pdf

Hi Richard,

We shipped 2 coolers today, the FedEx tracking is 804557917397 and 780047104871.

For ES107UF, "field filtered" should not be marked. The correction was made on the attached COC.

Thanks, Ann

Click here to report this email as spam.

130 140 CHAIN OF CUSTODY RECORD 01/01/14 Revision F 1200 SAMPLER(S): (PRINT) (3'002 N89 Time: Time: 8.812 [] 6817 [] 3617 [] (IV)10 REQUESTED ANALYSES 12114 X747\0209 \Big X747\0109 \Big sisteM SST P.O. NO.: MIC 07S8 TO 0YS8 CD 2HA9 Please check box or fill in blank as needed. Date: PCBs (8082) Pesticides (8081) Page\_ Date 800cs (8270) Prep (5035) 🗆 En Core 🗆 Terra Core Oxygenates (8260) **つつ021** CLIENT PROJECT NAME / NUMBER: ( hong OCs (8580) □ 81EX / MTBE □ 8260 □ WO # / LAB USE ONLY PROJECT CONTACT 24 1111 Poperty TPH C6-C36 C6-C44 Received by: (Signature/Affiliation) Received by: (Signature/Affiliation) Received by: (Signature/Affiliation) (5/2) DECTED (b)Hard (1928) <del>- ОЯЭ-Б-</del> (в)нчт**Ж** TELSO8-2 COL. UTUD Churu PESCIEMET-CUM AFCHERBETTICKED.CAM WE Calscience Environmental Laboratories, Inc. Field Filtered CHES. CLETZY LOG CODE X Preserved Χ LABORATORY CLIENT FAVI PO NYMEPIAI SCICNCE INTORNATIONA Other locations: Concord, San Luis Obispo, Houston, and Corpus Christi Unpreserved 7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494 X STANDARD S o S P S SNT. 9 σ For courier service / sample drop off information, contact sales@calscience.com or call us. REGULD (FULL) (AM MONITORING ANOTYTE 115) STATE: \*. 4 × 3 × 4 1. A CHOSE MATRIX TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"): ☐5 DAYS 080 120 のなら 1035 TIME 200 SAMPLING シェク 112711V HIZZIL 172712 ☐72 HR がんだし DATE 7 ☐48 HR ADDRESS: 3574 (ILMINU GLOBAL ID S 107 UT SAMPLE ID Relinquished by: (Signature) ☐24 HR Relinquished by: (Signature) Relinquished by: (Signature) 557RM SPECIAL INSTRUCTIONS: んさせ ハヘこと 2106 OITY: YALTUG ☐ COELT EDF SAME DAY LAB USE ONLY

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KAILUA: HI 967342532 UNITED STATES US

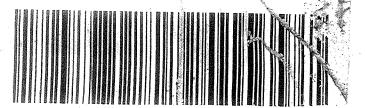
SHIP DATE: 22JUL14 ACTWGT: 48.0 LB CAD: /POS1501 DIMS: 24x13x13 IN

TO SAMPLE CONTROL CAL SCIENCE 7440 LINCOLN WAY

# **GARDEN GROVE CA 92841**

1 of 2 TRK# 0200 8045 5791 7397

WED - 23 JUL AA



354 ULUNIU ST STE 304

KAILUA, HI 967342532 UNITED STATES US

BILL SENCER

**TO SAMPLE CONTROL** CAL SCIENCE 7440 LINCOLN WAY

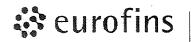
# GARDEN GROVE CA 92841 (714) 895-5494 REF:

STANDARD OVERNIGHT

WZ APVA

92841 CA-US SNA





# Calscience

WORK ORDER #: 14-07- ☐ 5 5 0

# SAMPLE RECEIPT FORM

Cooler \_\_\_ of \_2

CLIENT: Env. Science.	DATE: _	0//23	<u>/ 14</u>
TEMPERATURE: Thermometer ID: SC1 (Criteria: 0.0 °C – 6.0 °C, not froze	n except se	diment/tissu	e)
_	Blank	☐ Sample	
☐ Sample(s) outside temperature criteria (PM/APM contacted by:)		·	
☐ Sample(s) outside temperature criteria but received on ice/chilled on same d	ay of samp	lina.	
☐ Received at ambient temperature, placed on ice for transport by Co		91	
	onici.	Checked b	15
Ambient Temperature:   Air   Filter	,	Checked b	· ·
CUSTODY SEALS INTACT:			<u>,</u> ~
Cooler □ □ No (Not Intact) □ Not Present	□ N/A	Checked b	y: 15
☑ Sample □ □ No (Not Intact) □ Not Present		Checked b	y: <u>816</u>
			2.140
SAMPLE CONDITION:	Yes	No	N/A
Chain-Of-Custody (COC) document(s) received with samples		∐ 	
COC document(s) received complete			
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.			
□ No analysis requested. □ Not relinquished. □ No date/time relinquished.			
Sampler's name indicated on COC	•		
Sample container label(s) consistent with COC			
Sample container(s) intact and good condition	,		
Proper containers and sufficient volume for analyses requested			Ш
Analyses received within holding time			
Aqueous samples received within 15-minute holding time			
□ pH □ Residual Chlorine □ Dissolved Sulfides □ Dissolved Oxygen	•		
Proper preservation noted on COC or sample container	. 🖊		
☑ Unpreserved vials received for Volatiles analysis	/		
Volatile analysis container(s) free of headspace			
Tedlar bag(s) free of condensation  CONTAINER TYPE:			
Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve () □EnCore	s <sup>®</sup> □Terra	aCores <sup>®</sup> □_	
Aqueous: ØVOA ØVOAh □VOAna₂ □125AGB □125AGBh □125AGBp	Z1AGB	□1AGEna <sub>2</sub> [	□1AGB <b>s</b>
□500AGB ☑500AGJ □500AGJs □250AGB □250CGB □250CGBs	s □1PB	□1PBna □	]500PB
☐250PB ☑250PBnu☐125PB ☐125PBznna ☐100PJ ☐100PJna; ☐		O.	
Air: Tedlar® Canister Other: Trip Blank Lot#: 131007B  Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: El	— nvelope	d/Checked by Reviewed by Scanned by	: 778



# Calscience

WORK ORDER #: 14-07- [ 5 5 0

# SAMPLE RECEIPT FORM

Cooler  $\underline{\mathcal{L}}$  of  $\underline{\mathcal{L}}$ 

CLIENT: Env. Science.	DATE:	07/23	/14
TEMPERATURE: Thermometer ID: SC1 (Criteria: 0.0 °C – 6.0 °C, not froze	n except se	ediment/tissu	ıe)
Temperature 2 • °C - 0.3 °C (CF) = 2 • 5 °C	Blank	☐ Sampl	е
☐ Sample(s) outside temperature criteria (PM/APM contacted by:)			
☐ Sample(s) outside temperature criteria but received on ice/chilled on same o	day of samp	ling.	
☐ Received at ambient temperature, placed on ice for transport by Co			
Ambient Temperature: □ Air □ Filter		Checked b	v: 15
CUSTODY SEALS INTACT:			
☐ Cooler ☐ ☐ No (Not Intact) ☐ Not Present	□ N/A	Checked b	y: <u>ا ک</u>
☑ Sample □ □ No (Not Intact) □ Not Present		Checked b	y: <u>846</u>
SAMPLE CONDITION:	Yes	No	N/A
Chain-Of-Custody (COC) document(s) received with samples			
COC document(s) received complete	,		
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels	<i>E.</i>	L <sup>2</sup>	_
☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.			
Sampler's name indicated on COC	. 🗹		
Sample container label(s) consistent with COC			
Sample container(s) intact and good condition			
Proper containers and sufficient volume for analyses requested			
Analyses received within holding time	_		
Aqueous samples received within 15-minute holding time			
☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen	. 🗆		Z
Proper preservation noted on COC or sample container	. 🖊		
Unpreserved vials received for Volatiles analysis	· _		
Volatile analysis container(s) free of headspace	. 🗹		
Tedlar bag(s) free of condensation  CONTAINER TYPE:	. 🗆		9
Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve () □EnCore	es <sup>®</sup> □Ţerra	aCores <sup>®</sup> □_	
Aqueous: ∠VOA □VOAh □VOAna₂ □125AGB □125AGBh □125AGBp	⊿1AGB I	□1AGB <b>na</b> ₂ [	∃1AGB <b>s</b>
□500AGB Ø500AGJ □500AGJs □250AGB □250CGB □250CGBs	s □1PB	□1PB <b>na</b> □	]500PB
□250PB №250PBnw□125PB □125PB <b>znna</b> □100PJ □100PJ <b>na</b> ₂ □		П_	· · · · · · · · · · · · · · · · · · ·
Air: □Tedlar <sup>®</sup> □Canister <b>Other:</b> □ Trip Blank Lot#:			
Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: El Preservative: h: HCL n: HNO <sub>3</sub> na <sub>2</sub> :Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> na: NaOH p: H <sub>3</sub> PO <sub>4</sub> s: H <sub>2</sub> SO <sub>4</sub> u: Ultra-pure znna: ZnAc <sub>2</sub> +Na		Reviewed by: Scanned by	



# Calscience

Supplemental Report 2

The original report has been revised/corrected.



# **WORK ORDER NUMBER: 14-07-1550**

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

**Analytical Report For** 

Client: Environmental Science International, Inc.

Client Project Name: Red Hill 112066

**Attention:** Robert Chong

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Richard Vellas

Approved for release on 07/31/2014 by: Richard Villafania

**Project Manager** 



ResultLink > Email your PM >

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.



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#### **Work Order Narrative**

Work Order: 14-07-1550 Page 1 of 1

#### **Condition Upon Receipt:**

Samples were received under Chain-of-Custody (COC) on 07/23/14. They were assigned to Work Order 14-07-1550.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

#### **Holding Times:**

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

#### **Quality Control:**

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

#### **Additional Comments:**

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New\_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

#### **Subcontractor Information:**

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.



 Environmental Science International, Inc.
 Date Received:
 07/23/14

 354 Uluniu Street, Suite 304
 Work Order:
 14-07-1550

 Kailua, HI 96734-2500
 Preparation:
 EPA 3510C

 Method:
 EPA 8015B (M)

 Units:
 ug/L

Project: Red Hill LTM 112066	Page 1 of 1

Client Sample Number		Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES106		14-07-1550-1-H	07/22/14 11:20	Aqueous	GC 45	07/24/14	07/25/14 09:09	140724B11B
Comment(s):	<ul> <li>Results were evaluated to</li> <li>TPH as Diesel is quantified</li> </ul>	` , ,		the MDL (DI	_) but < RL (LC	Q), if found, a	re qualified with a	a "J" flag.
Parameter TPH as Diesel		Result 37	<u>DL</u> 11	<u>LOD</u> 12	<u>LC</u> 25		<u>DF</u> 1.00	<u>Qualifiers</u> HD
Surrogate n-Octacosane		<u>Rec. (%)</u> 91	Control Limi 51-141	<u>its</u> Quali	<u>fiers</u>			

ES107	14-0	7-1550-2-G	07/22/14 09:45	Aqueous	GC 45	07/24/14	07/25/14 09:28	140724B11E
Comment(s):	- Results were evaluated to the N	MDL (DL), conc	entrations >= to the	ne MDL (DI	_) but < RL	(LOQ), if found, a	are qualified with a	a "J" flag.
	- TPH as Diesel is quantified in tl	he carbon range	e C10-C28.					
<u>Parameter</u>		Result	<u>DL</u>	<u>LOD</u>		<u>LOQ</u>	<u>DF</u>	<b>Qualifiers</b>
TPH as Diesel		<12	11	12		25	1.00	U
<u>Surrogate</u>		Rec. (%)	Control Limit	ts Quali	<u>fiers</u>			
n-Octacosane		92	51-141					

ES108	14-07-1550-3-F	07/22/14 10:35	Aqueous	GC 45	07/24/14	07/25/14 09:47	140724B11B
Comment(s):	- Results were evaluated to the MDL (DL), o	oncentrations >=	to the MDL (DL	) but < RL (	LOQ), if found, a	re qualified with	h a "J" flag.
	- TPH as Diesel is quantified in the carbon re	ange C10-C28.					
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>		<u>LOQ</u>	<u>DF</u>	<b>Qualifiers</b>
TPH as Diesel	<12	11	12		25	1.00	U

SurrogateRec. (%)Control LimitsQualifiersn-Octacosane10051-141

Method Blank	099-15-516-159	N/A	Aqueous	GC 45	07/24/14	07/25/14 03:30	140724B11B
Comment(s):	- Results were evaluated to the MDL (DL), con	centrations >= to	the MDL (DL	_) but < RL (	LOQ), if found, a	are qualified with a	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>		<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
TPH as Diesel	<12	11	12		25	1.00	U
Surrogate	Rec. (%)	Control Lim	<u>its Qualit</u>	<u>fiers</u>			
n-Octacosane	90	51-141					



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

07/23/14 14-07-1550

Work Order: Preparation:

EPA 3005A Filt.

Method:

EPA 6020

Units:

ug/L

Project: Red Hill LTM 112066

Page 1 of 1

Client Sample N	lumber	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES106		14-07-1550-1-G	07/22/14 11:20	Aqueous	ICP/MS 04	07/23/14	07/24/14 19:53	140723L03D
Comment(s):	- Results were evaluated to	the MDL (DL), cond	centrations >= to	o the MDL (DL	) but < RL (LO	Q), if found, a	re qualified with a	a "J" flag.
<u>Parameter</u>		Result	<u>DL</u>	<u>LOD</u>	<u>LO</u>	<u>Q</u>	<u>DF</u>	<u>Qualifiers</u>
Lead		<0.200	0.0898	0.200	1.0	0	1.00	U

ES108	14-07-1550-3-G	07/22/14 10:35	Aqueous	ICP/MS 04	07/23/14	07/24/14 20:00	140723L03D
Comment(s):	- Results were evaluated to the MDL (DL), con	centrations >= t	o the MDL (DL	_) but < RL (LO	Q), if found, are	e qualified with	a "J" flag.
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOG	<u>2</u>	<u>DF</u>	<u>Qualifiers</u>
Lead	<0.200	0.0898	0.200	1.00	)	1.00	U

Method Blank	099-14-497-87	7 N/A	Aqueous	ICP/MS 04	07/23/14	07/24/14 19:24	140723L03D
Comment(s):	- Results were evaluated to the MDL (DL),	concentrations	s >= to the MDL (DL	) but < RL (LOQ	), if found, are	e qualified with a	"J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u> </u>	<u>DF</u>	<u>Qualifiers</u>
Lead	<0.200	0.08	98 0.200	1.00		1.00	U





Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received:

07/23/14 14-07-1550

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 1 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID	
ES106	14-07-1550-1-J	07/22/14 11:20	Aqueous	GC/MS AAA	07/28/14	07/29/14 17:29	140728L01	
Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.								
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOC	<u>)</u>	<u>DF</u>	<u>Qualifiers</u>	
Naphthalene	< 0.047	0.022	0.047	0.19	1	1.00	U	
2-Methylnaphthalene	< 0.047	0.025	0.047	0.19	1	1.00	U	
1-Methylnaphthalene	< 0.047	0.027	0.047	0.19	1	1.00	U	
Acenaphthylene	< 0.047	0.017	0.047	0.19	1	1.00	U	
Acenaphthene	< 0.047	0.020	0.047	0.19	)	1.00	U	
Fluorene	< 0.047	0.023	0.047	0.19	1	1.00	U	
Phenanthrene	< 0.047	0.029	0.047	0.19	)	1.00	U	
Anthracene	< 0.047	0.032	0.047	0.19	)	1.00	U	
Fluoranthene	< 0.047	0.026	0.047	0.19	)	1.00	U	
Pyrene	< 0.047	0.023	0.047	0.19	)	1.00	U	
Benzo (a) Anthracene	< 0.047	0.022	0.047	0.19	)	1.00	U	
Chrysene	< 0.047	0.018	0.047	0.19	)	1.00	U	
Benzo (k) Fluoranthene	< 0.047	0.022	0.047	0.19	)	1.00	U	
Benzo (b) Fluoranthene	< 0.047	0.024	0.047	0.19	)	1.00	U	
Benzo (a) Pyrene	< 0.047	0.034	0.047	0.19	)	1.00	U	
Indeno (1,2,3-c,d) Pyrene	< 0.047	0.021	0.047	0.19	)	1.00	U	
Dibenz (a,h) Anthracene	< 0.047	0.025	0.047	0.19	)	1.00	U	
Benzo (g,h,i) Perylene	<0.047	0.021	0.047	0.19	)	1.00	U	
Surrogate	Rec. (%)	Control Limi	its Qualifi	<u>iers</u>				
Nitrobenzene-d5	78	28-139						
2-Fluorobiphenyl	78	33-144						
p-Terphenyl-d14	78	23-160						



Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received:

07/23/14

Work Order: Preparation:

14-07-1550 EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 2 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID		
ES107	14-07-1550-2-H	07/22/14 09:45	Aqueous	GC/MS AAA	07/28/14	07/29/14 17:53	140728L01		
Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.									
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOC	<u>)</u>	<u>DF</u>	<u>Qualifiers</u>		
Naphthalene	<0.048	0.022	0.048	0.19	)	1.00	U		
2-Methylnaphthalene	<0.048	0.025	0.048	0.19	)	1.00	U		
1-Methylnaphthalene	<0.048	0.027	0.048	0.19	)	1.00	U		
Acenaphthylene	<0.048	0.017	0.048	0.19	)	1.00	U		
Acenaphthene	<0.048	0.020	0.048	0.19	)	1.00	U		
Fluorene	<0.048	0.023	0.048	0.19	)	1.00	U		
Phenanthrene	<0.048	0.029	0.048	0.19	)	1.00	U		
Anthracene	<0.048	0.033	0.048	0.19	)	1.00	U		
Fluoranthene	<0.048	0.026	0.048	0.19	)	1.00	U		
Pyrene	<0.048	0.024	0.048	0.19	)	1.00	U		
Benzo (a) Anthracene	<0.048	0.023	0.048	0.19	)	1.00	U		
Chrysene	<0.048	0.018	0.048	0.19	)	1.00	U		
Benzo (k) Fluoranthene	<0.048	0.022	0.048	0.19	)	1.00	U		
Benzo (b) Fluoranthene	<0.048	0.024	0.048	0.19	)	1.00	U		
Benzo (a) Pyrene	<0.048	0.035	0.048	0.19	)	1.00	U		
Indeno (1,2,3-c,d) Pyrene	<0.048	0.021	0.048	0.19	)	1.00	U		
Dibenz (a,h) Anthracene	<0.048	0.026	0.048	0.19	)	1.00	U		
Benzo (g,h,i) Perylene	<0.048	0.021	0.048	0.19	)	1.00	U		
Surrogate	Rec. (%)	Control Limi	its Qualifi	<u>iers</u>					
Nitrobenzene-d5	73	28-139							
2-Fluorobiphenyl	76	33-144							
p-Terphenyl-d14	71	23-160							



Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received: Work Order:

07/23/14 14-07-1550

Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 3 of 4

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID	
ES108	14-07-1550-3-J	07/22/14 10:35	Aqueous	GC/MS AAA	07/28/14	07/29/14 18:18	140728L01	
Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.								
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOC	<u>Q</u>	<u>DF</u>	<u>Qualifiers</u>	
Naphthalene	<0.049	0.023	0.049	0.20	)	1.00	U	
2-Methylnaphthalene	< 0.049	0.026	0.049	0.20	)	1.00	U	
1-Methylnaphthalene	<0.049	0.028	0.049	0.20	)	1.00	U	
Acenaphthylene	< 0.049	0.018	0.049	0.20	)	1.00	U	
Acenaphthene	< 0.049	0.020	0.049	0.20	)	1.00	U	
Fluorene	< 0.049	0.024	0.049	0.20	)	1.00	U	
Phenanthrene	< 0.049	0.030	0.049	0.20	)	1.00	U	
Anthracene	< 0.049	0.034	0.049	0.20	)	1.00	U	
Fluoranthene	< 0.049	0.027	0.049	0.20	)	1.00	U	
Pyrene	< 0.049	0.024	0.049	0.20	)	1.00	U	
Benzo (a) Anthracene	< 0.049	0.023	0.049	0.20	)	1.00	U	
Chrysene	< 0.049	0.019	0.049	0.20	)	1.00	U	
Benzo (k) Fluoranthene	< 0.049	0.023	0.049	0.20	)	1.00	U	
Benzo (b) Fluoranthene	< 0.049	0.024	0.049	0.20	)	1.00	U	
Benzo (a) Pyrene	< 0.049	0.036	0.049	0.20	)	1.00	U	
Indeno (1,2,3-c,d) Pyrene	< 0.049	0.022	0.049	0.20	)	1.00	U	
Dibenz (a,h) Anthracene	< 0.049	0.026	0.049	0.20	)	1.00	U	
Benzo (g,h,i) Perylene	<0.049	0.021	0.049	0.20	)	1.00	U	
Surrogate	Rec. (%)	Control Lim	nits Qualif	<u>iers</u>				
Nitrobenzene-d5	68	28-139						
2-Fluorobiphenyl	72	33-144						
p-Terphenyl-d14	64	23-160						



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

2-Fluorobiphenyl

p-Terphenyl-d14

Date Received: Work Order:

07/23/14 14-07-1550

Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L Page 4 of 4

Project: Red Hill LTM 112066

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID		
Method Blank	099-15-148-52	N/A	Aqueous	GC/MS AAA	07/28/14	07/29/14 13:03	140728L01		
Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.									
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOC	<u>)</u>	<u>DF</u>	<u>Qualifiers</u>		
Naphthalene	< 0.050	0.023	0.050	0.20	)	1.00	U		
2-Methylnaphthalene	< 0.050	0.026	0.050	0.20	)	1.00	U		
1-Methylnaphthalene	< 0.050	0.028	0.050	0.20	)	1.00	U		
Acenaphthylene	< 0.050	0.018	0.050	0.20	)	1.00	U		
Acenaphthene	< 0.050	0.021	0.050	0.20	)	1.00	U		
Fluorene	< 0.050	0.024	0.050	0.20	)	1.00	U		
Phenanthrene	< 0.050	0.031	0.050	0.20	)	1.00	U		
Anthracene	< 0.050	0.034	0.050	0.20	)	1.00	U		
Fluoranthene	< 0.050	0.027	0.050	0.20	)	1.00	U		
Pyrene	< 0.050	0.025	0.050	0.20	)	1.00	U		
Benzo (a) Anthracene	< 0.050	0.024	0.050	0.20	)	1.00	U		
Chrysene	< 0.050	0.019	0.050	0.20	)	1.00	U		
Benzo (k) Fluoranthene	< 0.050	0.023	0.050	0.20	)	1.00	U		
Benzo (b) Fluoranthene	< 0.050	0.025	0.050	0.20	)	1.00	U		
Benzo (a) Pyrene	< 0.050	0.036	0.050	0.20	)	1.00	U		
Indeno (1,2,3-c,d) Pyrene	< 0.050	0.022	0.050	0.20	)	1.00	U		
Dibenz (a,h) Anthracene	< 0.050	0.027	0.050	0.20	)	1.00	U		
Benzo (g,h,i) Perylene	<0.050	0.022	0.050	0.20	)	1.00	U		
Surrogate	Rec. (%)	Control Limi	<u>its</u> Qualifi	ers					
Nitrobenzene-d5	90	28-139							

33-144

23-160

86

87



Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received:

07/23/14 14-07-1550

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID	
ES106	14-07-1550-1-A	07/22/14 11:20	Aqueous	GC/MS OO	07/23/14	07/23/14 17:27	140723L017	
Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.								
<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	LOC	<u> </u>	<u>)F</u>	<b>Qualifiers</b>	
Acetone	<10	6.0	10	20	1	.00	U,ICH	
Benzene	<0.50	0.14	0.50	1.0	1	.00	U	
Bromodichloromethane	<0.50	0.21	0.50	5.0	1	.00	U	
Bromoform	<1.0	0.50	1.0	10	1	.00	U	
Bromomethane	<5.0	3.9	5.0	20	1	.00	U	
2-Butanone	<5.0	2.2	5.0	10	1	.00	U	
Carbon Tetrachloride	<0.50	0.23	0.50	1.0	1	.00	U	
Chlorobenzene	<0.50	0.17	0.50	5.0	1	.00	U	
Chloroethane	<5.0	2.3	5.0	10	1	.00	U	
Chloroform	<0.50	0.46	0.50	5.0	1	.00	U	
Chloromethane	<2.0	1.8	2.0	10	1	.00	U,IJ	
Dibromochloromethane	<0.50	0.25	0.50	1.0	1	.00	U	
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1	.00	U	
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1	.00	U	
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1	.00	U	
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1	.00	U	
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0	1	.00	U	
1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1	.00	U	
1,2-Dichloroethane	<0.50	0.24	0.50	1.0	1	.00	U	
1,1-Dichloroethene	<0.50	0.43	0.50	1.0	1	.00	U	
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0	1	.00	U	
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0	1	.00	U	
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1	.00	U	
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1	.00	U	
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1	.00	U	
Ethylbenzene	<0.50	0.14	0.50	1.0	1	.00	U	
Methylene Chloride	<1.0	0.64	1.0	5.0	1	.00	U	
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1	.00	U	
Styrene	<0.50	0.17	0.50	1.0	1	.00	U	
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1	.00	U	
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0	1	.00	U	
Tetrachloroethene	<0.50	0.39	0.50	5.0	1	.00	U	
Toluene	<0.50	0.24	0.50	1.0		.00	U	
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0		.00	U	
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0		.00	U	
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0		.00	U	



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Date Received: Work Order: Preparation: Method:

07/23/14 14-07-1550 EPA 5030C GC/MS / EPA 8260B

Units:

ug/L Page 2 of 10

Project: Red Hill LTM 112066

<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U

Surrogate	Rec. (%)	Control Limits	<b>Qualifiers</b>
Dibromofluoromethane	97	80-126	
1,2-Dichloroethane-d4	90	80-134	
Toluene-d8	98	80-120	
Toluene-d8-TPPH	97	88-112	
1,4-Bromofluorobenzene	94	80-120	





Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

07/23/14 14-07-1550

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES107	14-07-1550-2-A	07/22/14 09:45	Aqueous	GC/MS OO	07/23/14	07/23/14 21:31	140723L017
Comment(s): - Results were evaluated to	the MDL (DL), con	centrations >= t	o the MDL (DL	) but < RL (LOC	), if found, are o	qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>DF</u>	=	<u>Qualifiers</u>
Acetone	<10	6.0	10	20	1.0	00	U,ICH
Benzene	< 0.50	0.14	0.50	1.0	1.0	00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0	1.0	00	U
Bromoform	<1.0	0.50	1.0	10	1.0	00	U
Bromomethane	<5.0	3.9	5.0	20	1.0	00	U
2-Butanone	<5.0	2.2	5.0	10	1.0	00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0	1.0	00	U
Chlorobenzene	<0.50	0.17	0.50	5.0	1.0	00	U
Chloroethane	<5.0	2.3	5.0	10	1.0	00	U
Chloroform	<0.50	0.46	0.50	5.0	1.0	00	U
Chloromethane	<2.0	1.8	2.0	10	1.0	00	U,IJ
Dibromochloromethane	<0.50	0.25	0.50	1.0	1.0	00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1.0	00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0	1.0	00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0	1.0	00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0	1.0	00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0	1.0	00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0	1.0	00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0	1.0	00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0	1.0	00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0	1.0	00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0	1.0	00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0	1.0	00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.0	00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0	1.0	00	U
Ethylbenzene	<0.50	0.14	0.50	1.0	1.0	00	U
Methylene Chloride	<1.0	0.64	1.0	5.0	1.0	00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1.0	00	U
Styrene	<0.50	0.17	0.50	1.0	1.0	00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1.0		U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0	1.0		U
Tetrachloroethene	<0.50	0.39	0.50	5.0	1.0		U
Toluene	<0.50	0.24	0.50	1.0	1.0		U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1.0		U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0	1.0		U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0	1.0		U



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Date Received: Work Order: Preparation: Method:

<u>LOD</u>

<u>LOQ</u>

1.0

1.0 5.0

1.0 10

1.0

1.0

50

<u>DF</u>

1.00 1.00

1.00

1.00

1.00

1.00

1.00 1.00

07/23/14 14-07-1550 EPA 5030C GC/MS / EPA 8260B

Units:

<u>DL</u>

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Qualifiers U

U

U

U

U

U U

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ug/L

Project: Red Hill LTM 11206	6
<u>Parameter</u>	

1,1,2-Trichloroethane	<0.50	0.38	0.50
Trichloroethene	<0.50	0.37	0.50
1,2,3-Trichloropropane	<1.0	0.64	1.0
Vinyl Chloride	<0.50	0.30	0.50
p/m-Xylene	<1.0	0.30	1.0
o-Xylene	<0.50	0.23	0.50
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50
Gasoline Range Organics	<30	26	30
Surrogate	Rec. (%)	Control Limits	<u>Qualifiers</u>
Dibromofluoromethane	98	80-126	
		00 .20	
1,2-Dichloroethane-d4	94	80-134	
1,2-Dichloroethane-d4 Toluene-d8	94 101		
	_	80-134	

Result





Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

07/23/14 14-07-1550

Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

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Project: Red Hill LTM 112066

Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES108	14-07-1550-3-A	07/22/14 10:35	Aqueous	GC/MS OO	07/23/14	07/23/14 21:58	140723L017
Comment(s): - Results were evaluated	to the MDL (DL), con	centrations >= t	o the MDL (DI	_) but < RL (LOC	)), if found, ar	re qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOG	<u>)</u>	<u>DF</u>	<b>Qualifiers</b>
Acetone	<10	6.0	10	20		1.00	U,ICH
Benzene	<0.50	0.14	0.50	1.0		1.00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0		1.00	U
Bromoform	<1.0	0.50	1.0	10		1.00	U
Bromomethane	<5.0	3.9	5.0	20		1.00	U
2-Butanone	<5.0	2.2	5.0	10		1.00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0		1.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0		1.00	U
Chloroethane	<5.0	2.3	5.0	10		1.00	U
Chloroform	<0.50	0.46	0.50	5.0		1.00	U
Chloromethane	<2.0	1.8	2.0	10		1.00	U,IJ
Dibromochloromethane	<0.50	0.25	0.50	1.0		1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10		1.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0		1.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0		1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0		1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0		1.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0		1.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0		1.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0		1.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0		1.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0		1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0		1.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0		1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0		1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10		1.00	U
Styrene	<0.50	0.17	0.50	1.0		1.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0		1.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0		1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0		1.00	U
Toluene	<0.50	0.24	0.50	1.0		1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0		1.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0		1.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0		1.00	U



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Date Received: Work Order: Preparation: Method:

07/23/14 14-07-1550 EPA 5030C GC/MS / EPA 8260B

Units:

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Project: Red Hill LTM 112066

<u>Parameter</u>	<u>Result</u>	<u>DL</u>	<u>LOD</u>	<u>LOQ</u>	<u>DF</u>	<b>Qualifiers</b>
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U
Surrogate	Rec. (%)	Control Limits	<u>Qualifiers</u>			

<u>Surrogate</u>	Rec. (%)	Control Limits
Dibromofluoromethane	98	80-126
1,2-Dichloroethane-d4	92	80-134
Toluene-d8	100	80-120
Toluene-d8-TPPH	99	88-112
1,4-Bromofluorobenzene	94	80-120





Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received: Work Order:

14-07-1550

07/23/14

Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
ES TRIP	14-07-1550-4-A	07/22/14 08:00	Aqueous	GC/MS OO	07/23/14	07/23/14 21:03	140723L017
Comment(s): - Results were evaluated t	o the MDL (DL), con	centrations >= t	o the MDL (DL	but < RL (LOC	Q), if found, are q	ualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	LOD	LOC	<u>DF</u>	<u>.</u>	<u>Qualifiers</u>
Acetone	<10	6.0	10	20	1.0	00	U,ICH
Benzene	<0.50	0.14	0.50	1.0	1.0	00	U
Bromodichloromethane	< 0.50	0.21	0.50	5.0	1.0	00	U
Bromoform	<1.0	0.50	1.0	10	1.0	00	U
Bromomethane	<5.0	3.9	5.0	20	1.0	00	U
2-Butanone	<5.0	2.2	5.0	10	1.0	00	U
Carbon Tetrachloride	< 0.50	0.23	0.50	1.0	1.0	00	U
Chlorobenzene	< 0.50	0.17	0.50	5.0	1.0	00	U
Chloroethane	<5.0	2.3	5.0	10	1.0	00	U
Chloroform	< 0.50	0.46	0.50	5.0	1.0	00	U
Chloromethane	<2.0	1.8	2.0	10	1.0	00	U,IJ
Dibromochloromethane	< 0.50	0.25	0.50	1.0	1.0	00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10	1.0	00	U
1,2-Dibromoethane	< 0.50	0.36	0.50	1.0	1.0	00	U
1,2-Dichlorobenzene	< 0.50	0.46	0.50	1.0	1.0	00	U
1,3-Dichlorobenzene	< 0.50	0.40	0.50	1.0	1.0	00	U
1,4-Dichlorobenzene	< 0.50	0.43	0.50	1.0	1.0	00	U
1,1-Dichloroethane	< 0.50	0.28	0.50	5.0	1.0	00	U
1,2-Dichloroethane	< 0.50	0.24	0.50	1.0	1.0	00	U
1,1-Dichloroethene	< 0.50	0.43	0.50	1.0	1.0	00	U
c-1,2-Dichloroethene	< 0.50	0.48	0.50	1.0	1.0	00	U
t-1,2-Dichloroethene	< 0.50	0.37	0.50	1.0	1.0	00	U
1,2-Dichloropropane	< 0.50	0.42	0.50	5.0	1.0	00	U
c-1,3-Dichloropropene	< 0.50	0.25	0.50	1.0	1.0	00	U
t-1,3-Dichloropropene	< 0.50	0.25	0.50	1.0	1.0	00	U
Ethylbenzene	< 0.50	0.14	0.50	1.0	1.0	00	U
Methylene Chloride	<1.0	0.64	1.0	5.0	1.0	00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10	1.0	00	U
Styrene	<0.50	0.17	0.50	1.0	1.0	00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0	1.0	00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0	1.0	00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0	1.0	00	U
Toluene	<0.50	0.24	0.50	1.0	1.0	00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0	1.0	00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0	1.0	00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0	1.0	00	U



Project: Red Hill LTM 112066

# **Analytical Report**

Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Date Received: Work Order: Preparation: Method:

07/23/14 14-07-1550 EPA 5030C GC/MS / EPA 8260B

> Qualifiers U

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U

Units:

ug/L Page 8 of 10

<u>DF</u>

1.00

1.00

1.00

1.00

1.00

1.00

1.00 1.00

<u>LOQ</u>

1.0

1.0

5.0

1.0 10

1.0

1.0

50

<u>Result</u>	<u>DL</u>	<u>LOD</u>
<0.50	0.38	0.50
<0.50	0.37	0.50
<1.0	0.64	1.0
<0.50	0.30	0.50
<1.0	0.30	1.0
<0.50	0.23	0.50
<0.50	0.31	0.50
<30	26	30
Rec. (%)	Control Limits	<b>Qualifiers</b>
00	00.400	
98	80-126	
98	80-126 80-134	
93	80-134	
	<0.50 <0.50 <1.0 <0.50 <1.0 <0.50 <0.50 <0.50 <30	<ul> <li>&lt;0.50</li> <li>0.38</li> <li>&lt;0.50</li> <li>0.37</li> <li>&lt;1.0</li> <li>0.64</li> <li>&lt;0.50</li> <li>0.30</li> <li>&lt;1.0</li> <li>0.30</li> <li>&lt;0.50</li> <li>0.23</li> <li>&lt;0.50</li> <li>0.31</li> <li>&lt;30</li> <li>26</li> <li>Rec. (%)</li> <li>Control Limits</li> </ul>



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

07/23/14 14-07-1550

Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

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Client Sample Number	Lab Sample Number	Date/Time Collected	Matrix	Instrument	Date Prepared	Date/Time Analyzed	QC Batch ID
Method Blank	099-13-057-58	N/A	Aqueous	GC/MS OO	07/23/14	07/23/14 16:50	140723L017
Comment(s): - Results were evaluated t	to the MDL (DL), cor	ncentrations >= t	o the MDL (DI	_) but < RL (LOC	Q), if found, are	e qualified with	a "J" flag.
<u>Parameter</u>	Result	<u>DL</u>	<u>LOD</u>	LOC	<u>2</u>	<u>DF</u>	<b>Qualifiers</b>
Acetone	<10	6.0	10	20		1.00	U
Benzene	<0.50	0.14	0.50	1.0		1.00	U
Bromodichloromethane	<0.50	0.21	0.50	5.0		1.00	U
Bromoform	<1.0	0.50	1.0	10		1.00	U
Bromomethane	<5.0	3.9	5.0	20		1.00	U
2-Butanone	<5.0	2.2	5.0	10		1.00	U
Carbon Tetrachloride	<0.50	0.23	0.50	1.0		1.00	U
Chlorobenzene	<0.50	0.17	0.50	5.0		1.00	U
Chloroethane	<5.0	2.3	5.0	10		1.00	U
Chloroform	<0.50	0.46	0.50	5.0		1.00	U
Chloromethane	<2.0	1.8	2.0	10		1.00	U
Dibromochloromethane	<0.50	0.25	0.50	1.0		1.00	U
1,2-Dibromo-3-Chloropropane	<2.0	1.2	2.0	10		1.00	U
1,2-Dibromoethane	<0.50	0.36	0.50	1.0		1.00	U
1,2-Dichlorobenzene	<0.50	0.46	0.50	1.0		1.00	U
1,3-Dichlorobenzene	<0.50	0.40	0.50	1.0		1.00	U
1,4-Dichlorobenzene	<0.50	0.43	0.50	1.0		1.00	U
1,1-Dichloroethane	<0.50	0.28	0.50	5.0		1.00	U
1,2-Dichloroethane	<0.50	0.24	0.50	1.0		1.00	U
1,1-Dichloroethene	<0.50	0.43	0.50	1.0		1.00	U
c-1,2-Dichloroethene	<0.50	0.48	0.50	1.0		1.00	U
t-1,2-Dichloroethene	<0.50	0.37	0.50	1.0		1.00	U
1,2-Dichloropropane	<0.50	0.42	0.50	5.0		1.00	U
c-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
t-1,3-Dichloropropene	<0.50	0.25	0.50	1.0		1.00	U
Ethylbenzene	<0.50	0.14	0.50	1.0		1.00	U
Methylene Chloride	<1.0	0.64	1.0	5.0		1.00	U
4-Methyl-2-Pentanone	<5.0	4.4	5.0	10		1.00	U
Styrene	<0.50	0.17	0.50	1.0		1.00	U
1,1,1,2-Tetrachloroethane	<0.50	0.40	0.50	1.0		1.00	U
1,1,2,2-Tetrachloroethane	<0.50	0.41	0.50	1.0		1.00	U
Tetrachloroethene	<0.50	0.39	0.50	5.0		1.00	U
Toluene	<0.50	0.24	0.50	1.0		1.00	U
1,2,4-Trichlorobenzene	<1.0	0.50	1.0	5.0		1.00	U
1,1,1-Trichloroethane	<0.50	0.30	0.50	5.0		1.00	U
Hexachloro-1,3-Butadiene	<0.50	0.32	0.50	1.0		1.00	U



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

1,4-Bromofluorobenzene

Project: Red Hill LTM 112066

Date Received: Work Order: Preparation:

Method: Units: 07/23/14 14-07-1550 EPA 5030C

GC/MS / EPA 8260B

ug/L

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<u>Parameter</u>	Result	<u>DL</u>	LOD	<u>LOQ</u>	<u>DF</u>	<u>Qualifiers</u>
1,1,2-Trichloroethane	<0.50	0.38	0.50	1.0	1.00	U
Trichloroethene	<0.50	0.37	0.50	1.0	1.00	U
1,2,3-Trichloropropane	<1.0	0.64	1.0	5.0	1.00	U
Vinyl Chloride	<0.50	0.30	0.50	1.0	1.00	U
p/m-Xylene	<1.0	0.30	1.0	10	1.00	U
o-Xylene	<0.50	0.23	0.50	1.0	1.00	U
Methyl-t-Butyl Ether (MTBE)	<0.50	0.31	0.50	1.0	1.00	U
Gasoline Range Organics	<30	26	30	50	1.00	U
<u>Surrogate</u>	Rec. (%)	Control Limits	<u>Qualifiers</u>			
Dibromofluoromethane	97	80-126				
1,2-Dichloroethane-d4	90	80-134				
Toluene-d8	99	80-120				
Toluene-d8-TPPH	98	88-112				

80-120

95



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Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

Date Received:

07/23/14

Work Order:

14-07-1550

EPA 3005A Filt.

Method:

EPA 6020

Project: Red Hill LTM 112066 Page 1 of 4

Quality Control Sample ID	Type		Matrix	Ins	strument	Date Prepared	Date Ana	lyzed	MS/MSD Ba	tch Number
14-07-1466-2	Sample		Aqueous	s IC	P/MS 04	07/23/14	07/24/14	19:48	140723S03	
14-07-1466-2	Matrix Spike		Aqueous	s IC	P/MS 04	07/23/14	07/24/14	19:39	140723S03	
14-07-1466-2	Matrix Spike I	Duplicate	Aqueous	s IC	P/MS 04	07/23/14	07/24/14	19:41	140723S03	
Parameter	Sample Conc.	<u>Spike</u> Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Lead	ND	100.0	105.9	106	112.8	113	80-120	6	0-20	



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Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received: Work Order:

07/23/14 14-07-1550

Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

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Quality Control Sample ID	Туре		Matrix		nstrument	Date Prepared	Date Ana	lyzed	MS/MSD Ba	atch Number
14-07-1466-2	Sample		Aqueou	ıs (	GC/MS AAA	07/28/14	07/29/14	21:31	140728S01	
14-07-1466-2	Matrix Spike		Aqueou	ıs	GC/MS AAA	07/28/14	07/29/14	13:52	140728S01	
14-07-1466-2	Matrix Spike	Duplicate	Aqueou	ıs (	GC/MS AAA	07/28/14	07/29/14	14:16	140728S01	
<u>Parameter</u>	<u>Sample</u> <u>Conc.</u>	<u>Spike</u> <u>Added</u>	MS Conc.	MS %Red	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Naphthalene	70.83	2.000	59.97	0	64.80	0	21-133	8	0-25	3
2-Methylnaphthalene	19.64	2.000	21.30	83	18.58	0	21-140	14	0-25	3
1-Methylnaphthalene	25.19	2.000	25.26	4	24.73	0	20-140	2	0-25	3
Acenaphthylene	ND	2.000	1.370	68	1.495	75	33-145	9	0-25	
Acenaphthene	0.5154	2.000	1.776	63	1.936	71	49-121	9	0-25	
Fluorene	0.2363	2.000	1.536	65	1.689	73	59-121	10	0-25	
Phenanthrene	ND	2.000	1.313	66	1.517	76	54-120	14	0-25	
Anthracene	ND	2.000	1.421	71	1.551	78	27-133	9	0-25	
Fluoranthene	ND	2.000	1.282	64	1.431	72	26-137	11	0-25	
Pyrene	ND	2.000	1.192	60	1.344	67	18-168	12	0-25	
Benzo (a) Anthracene	ND	2.000	1.252	63	1.416	71	33-143	12	0-25	
Chrysene	ND	2.000	1.298	65	1.459	73	17-168	12	0-25	
Benzo (k) Fluoranthene	ND	2.000	1.137	57	1.284	64	24-159	12	0-25	
Benzo (b) Fluoranthene	ND	2.000	1.117	56	1.282	64	24-159	14	0-25	
Benzo (a) Pyrene	ND	2.000	1.181	59	1.348	67	17-163	13	0-25	
Indeno (1,2,3-c,d) Pyrene	ND	2.000	1.190	60	1.329	66	10-171	11	0-25	
Dibenz (a,h) Anthracene	ND	2.000	1.195	60	1.342	67	10-219	12	0-25	
Benzo (g,h,i) Perylene	ND	2.000	1.252	63	1.389	69	10-227	10	0-25	



Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

EPA 5030C

Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066	Page 3 of 4
1 TOJCCL TCG TIIII ETW T 12000	i age o or <del>-</del>

Quality Control Sample ID	Туре		Matrix	Ir	strument	Date Prepared	Date Ana	lyzed	MS/MSD Bat	tch Number
ES106	Sample		Aqueous	G	C/MS OO	07/23/14	07/23/14	17:27	140723S011	
ES106	Matrix Spike		Aqueous	G	C/MS OO	07/23/14	07/23/14	17:54	140723S011	
ES106	Matrix Spike	Duplicate	Aqueous	G	C/MS OO	07/23/14	07/23/14	18:21	140723S011	
<u>Parameter</u>	Sample Conc.	Spike Added	MS Conc.	MS %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
Acetone	ND	50.00	78.70	157	77.59	155	40-140	1	0-20	3
Benzene	ND	50.00	48.76	98	48.22	96	80-120	1	0-20	
Bromodichloromethane	ND	50.00	50.15	100	49.50	99	75-120	1	0-20	
Bromoform	ND	50.00	49.49	99	49.74	99	70-130	0	0-20	
Bromomethane	ND	50.00	48.54	97	44.16	88	30-145	9	0-20	
2-Butanone	ND	50.00	54.78	110	56.44	113	30-150	3	0-20	
Carbon Tetrachloride	ND	50.00	46.25	93	46.61	93	65-140	1	0-20	
Chlorobenzene	ND	50.00	50.80	102	49.94	100	80-120	2	0-20	
Chloroethane	ND	50.00	45.52	91	45.22	90	60-135	1	0-20	
Chloroform	ND	50.00	47.61	95	48.02	96	65-135	1	0-20	
Chloromethane	ND	50.00	41.38	83	41.46	83	40-125	0	0-20	
Dibromochloromethane	ND	50.00	50.97	102	51.09	102	60-135	0	0-20	
1,2-Dibromo-3-Chloropropane	ND	50.00	44.26	89	45.26	91	50-130	2	0-20	
1,2-Dibromoethane	ND	50.00	47.58	95	48.04	96	80-120	1	0-20	
1,2-Dichlorobenzene	ND	50.00	50.78	102	50.89	102	70-120	0	0-20	
1,3-Dichlorobenzene	ND	50.00	51.60	103	51.41	103	75-125	0	0-20	
1,4-Dichlorobenzene	ND	50.00	49.19	98	48.80	98	75-125	1	0-20	
1,1-Dichloroethane	ND	50.00	46.01	92	46.85	94	70-135	2	0-20	
1,2-Dichloroethane	ND	50.00	46.31	93	45.80	92	70-130	1	0-20	
1,1-Dichloroethene	ND	50.00	46.79	94	47.20	94	70-130	1	0-20	
c-1,2-Dichloroethene	ND	50.00	52.45	105	52.99	106	70-125	1	0-20	
t-1,2-Dichloroethene	ND	50.00	49.97	100	50.27	101	60-140	1	0-20	
1,2-Dichloropropane	ND	50.00	49.22	98	48.61	97	75-125	1	0-20	
c-1,3-Dichloropropene	ND	50.00	52.35	105	51.96	104	70-130	1	0-20	
t-1,3-Dichloropropene	ND	50.00	51.06	102	50.54	101	55-140	1	0-20	
Ethylbenzene	ND	50.00	49.48	99	49.20	98	75-125	1	0-20	
Methylene Chloride	ND	50.00	50.57	101	50.66	101	55-140	0	0-20	
4-Methyl-2-Pentanone	ND	50.00	48.94	98	50.04	100	60-135	2	0-20	
Styrene	ND	50.00	51.02	102	50.70	101	65-135	1	0-20	
1,1,1,2-Tetrachloroethane	ND	50.00	47.91	96	47.87	96	80-130	0	0-20	
1,1,2,2-Tetrachloroethane	ND	50.00	52.31	105	53.33	107	65-130	2	0-20	
Tetrachloroethene	ND	50.00	44.59	89	43.31	87	45-150	3	0-20	
Toluene	ND	50.00	48.57	97	48.18	96	75-120	1	0-20	
1,2,4-Trichlorobenzene	ND	50.00	49.23	98	49.82	100	65-135	1	0-20	
1,1,1-Trichloroethane	ND	50.00	45.95	92	46.57	93	65-130	1	0-20	



Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

Date Received:

07/23/14

Work Order:

14-07-1550

EPA 5030C

Project: Red Hill LTM 112066 Page 4 of 4

<u>Parameter</u>	Sample Conc.	<u>Spike</u> <u>Added</u>	MS Conc.	<u>MS</u> %Rec.	MSD Conc.	MSD %Rec.	%Rec. CL	<u>RPD</u>	RPD CL	<u>Qualifiers</u>
Hexachloro-1,3-Butadiene	ND	50.00	46.15	92	46.43	93	50-140	1	0-20	
1,1,2-Trichloroethane	ND	50.00	50.04	100	49.56	99	75-125	1	0-20	
Trichloroethene	ND	50.00	48.06	96	47.51	95	70-125	1	0-20	
1,2,3-Trichloropropane	ND	50.00	47.27	95	47.63	95	75-125	1	0-20	
Vinyl Chloride	ND	50.00	45.23	90	44.53	89	50-145	2	0-20	
p/m-Xylene	ND	100.0	98.81	99	97.46	97	75-130	1	0-20	
o-Xylene	ND	50.00	51.41	103	51.11	102	80-120	1	0-20	
Methyl-t-Butyl Ether (MTBE)	ND	50.00	45.34	91	46.70	93	65-125	3	0-20	



# **Quality Control - PDS**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

Date Received:

07/23/14

Work Order:

14-07-1550

EPA 3005A Filt.

Method:

EPA 6020

Project: Red Hill LTM 112066 Page 1 of 1

Quality Control Sample ID	Туре	N	/latrix	Instrument	Date Prepared	Date Analyzed	PDS/PDSD Batch Number
14-07-1466-2	Sample	Δ.	Aqueous	ICP/MS 04	07/23/14 00:00	07/24/14 19:48	140723S03
14-07-1466-2	PDS	A	Aqueous	ICP/MS 04	07/23/14 00:00	07/24/14 19:43	140723S03
<u>Parameter</u>		Sample Conc.	Spike Added	PDS Conc.	PDS %Re	ec. %Rec. C	<u>Qualifiers</u>
Lead		ND	100.0	102.3	102	75-125	





# **Quality Control - LCS/LCSD**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

Project: Red Hill LTM 112066

Date Received:

07/23/14

Work Order:

14-07-1550

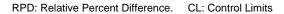
Preparation:

EPA 3510C

Method:

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Quality Control Sample ID	Туре	Mat	rix	Instrument	Date Pre	pared Date	e Analyzed	LCS/LCSD Ba	atch Number
099-15-516-159	LCS	Aqı	ueous	GC 45	07/24/14	07/2	5/14 03:49	140724B11B	
099-15-516-159	LCSD	Aqı	ieous	GC 45	07/24/14	07/2	5/14 04:08	140724B11B	
Parameter	Spike Added	LCS Conc.	LCS %Rec.	LCSD Conc.	LCSD %Rec.	%Rec. CL	RPD	RPD CL	Qualifiers
TPH as Diesel	2000	2008	100	2003	100	60-132	0	0-11	





# **Quality Control - LCS**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

Date Received:

07/23/14

Work Order:

14-07-1550

EPA 3005A Filt.

Method:

EPA 6020

Project: Red Hill LTM 112066 Page 2 of 5

Quality Control Sample ID	Туре	Matrix	Instrument	Date Prepared	Date Analyzed	LCS Batch Number
099-14-497-87	LCS	Aqueous	ICP/MS 04	07/23/14	07/24/14 19:32	140723L03D
Parameter		Spike Added	Conc. Recovere	ed LCS %Re	ec. %Rec	. CL Qualifiers
Lead		100.0	91.75	92	80-12	0



# **Quality Control - LCS**

Environmental Science International, Inc.

Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received: Work Order:

07/23/14 14-07-1550

354 Uluniu Street, Suite 304

Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

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Quality Control Sample ID	Туре	Matrix	Instrument	Date Prepared D	ate Analyzed LCS Ba	tch Number
099-15-148-52	LCS	Aqueous	GC/MS AAA	07/28/14 0	7/29/14 13:28 140728	L01
<u>Parameter</u>		Spike Added	Conc. Recovere	ed LCS %Rec.	%Rec. CL	<b>Qualifiers</b>
Naphthalene		2.000	1.361	68	21-133	
2-Methylnaphthalene		2.000	1.230	61	21-140	
1-Methylnaphthalene		2.000	1.226	61	20-140	
Acenaphthylene		2.000	1.167	58	33-145	
Acenaphthene		2.000	1.271	64	55-121	
Fluorene		2.000	1.315	66	59-121	
Phenanthrene		2.000	1.379	69	54-120	
Anthracene		2.000	1.393	70	27-133	
Fluoranthene		2.000	1.385	69	26-137	
Pyrene		2.000	1.333	67	45-129	
Benzo (a) Anthracene		2.000	1.343	67	33-143	
Chrysene		2.000	1.447	72	17-168	
Benzo (k) Fluoranthene		2.000	1.265	63	24-159	
Benzo (b) Fluoranthene		2.000	1.294	65	24-159	
Benzo (a) Pyrene		2.000	1.273	64	17-163	
Indeno (1,2,3-c,d) Pyrene		2.000	1.408	70	25-175	
Dibenz (a,h) Anthracene		2.000	1.350	67	25-175	
Benzo (g,h,i) Perylene		2.000	1.500	75	25-157	



# **Quality Control - LCS/LCSD**

Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

07/23/14 14-07-1550

Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Project: Red Hill LTM 1120	66					Page 4 of 5
Quality Control Sample ID	Type	Matrix	Instrument	Date Prepared	Date Analyzed	LCS/LCSD Batch Number

Quality Control Sample ID	Туре	Mat	rix	Instrument	Date Pre	pared	Date	Analyzed	LCS/LCSD Ba	atch Number
099-13-057-58	LCS	Aqı	ueous	GC/MS OO	07/23/14		07/23	/14 15:27	140723L017	
099-13-057-58	LCSD	Aqı	ieous	GC/MS OO	07/23/14		07/23	/14 15:54	140723L017	
Parameter	Spike Added	LCS Conc.	LCS %Rec.	LCSD Conc.	LCSD %Rec.	%Red	. CL	RPD	RPD CL	Qualifiers
Acetone	50.00	76.54	153	N/A	N/A	40-14	0	N/A	0-20	Χ
Benzene	50.00	48.59	97	N/A	N/A	80-12	0	N/A	0-20	
Bromodichloromethane	50.00	50.08	100	N/A	N/A	75-12	0	N/A	0-20	
Bromoform	50.00	50.09	100	N/A	N/A	70-13	0	N/A	0-20	
Bromomethane	50.00	46.74	93	N/A	N/A	30-14	5	N/A	0-20	
2-Butanone	50.00	56.87	114	N/A	N/A	30-15	0	N/A	0-20	
Carbon Tetrachloride	50.00	46.09	92	N/A	N/A	65-14	0	N/A	0-20	
Chlorobenzene	50.00	51.12	102	N/A	N/A	80-12	0	N/A	0-20	
Chloroethane	50.00	44.84	90	N/A	N/A	60-13	5	N/A	0-20	
Chloroform	50.00	48.48	97	N/A	N/A	65-13	5	N/A	0-20	
Chloromethane	50.00	41.19	82	N/A	N/A	40-12	5	N/A	0-20	
Dibromochloromethane	50.00	52.03	104	N/A	N/A	60-13	5	N/A	0-20	
1,2-Dibromo-3-Chloropropane	50.00	45.40	91	N/A	N/A	50-13	0	N/A	0-20	
1,2-Dibromoethane	50.00	48.92	98	N/A	N/A	80-12	0	N/A	0-20	
1,2-Dichlorobenzene	50.00	51.03	102	N/A	N/A	70-12	0	N/A	0-20	
1,3-Dichlorobenzene	50.00	51.52	103	N/A	N/A	75-12	5	N/A	0-20	
1,4-Dichlorobenzene	50.00	48.28	97	N/A	N/A	75-12	5	N/A	0-20	
1,1-Dichloroethane	50.00	47.01	94	N/A	N/A	70-13	5	N/A	0-20	
1,2-Dichloroethane	50.00	46.72	93	N/A	N/A	70-13	0	N/A	0-20	
1,1-Dichloroethene	50.00	47.70	95	N/A	N/A	70-13	0	N/A	0-20	
c-1,2-Dichloroethene	50.00	53.05	106	N/A	N/A	70-12	5	N/A	0-20	
t-1,2-Dichloroethene	50.00	51.14	102	N/A	N/A	60-14	0	N/A	0-20	
1,2-Dichloropropane	50.00	49.85	100	N/A	N/A	75-12	5	N/A	0-20	
c-1,3-Dichloropropene	50.00	52.80	106	N/A	N/A	70-13	0	N/A	0-20	
t-1,3-Dichloropropene	50.00	51.91	104	N/A	N/A	55-14	0	N/A	0-20	
Ethylbenzene	50.00	49.38	99	N/A	N/A	75-12	5	N/A	0-20	
Methylene Chloride	50.00	51.00	102	N/A	N/A	55-14	0	N/A	0-20	
4-Methyl-2-Pentanone	50.00	50.30	101	N/A	N/A	60-13	5	N/A	0-20	
Styrene	50.00	51.16	102	N/A	N/A	65-13	5	N/A	0-20	
1,1,1,2-Tetrachloroethane	50.00	48.31	97	N/A	N/A	80-13	0	N/A	0-20	
1,1,2,2-Tetrachloroethane	50.00	50.35	101	N/A	N/A	65-13	0	N/A	0-20	
Tetrachloroethene	50.00	49.33	99	N/A	N/A	45-15	0	N/A	0-20	
Toluene	50.00	48.73	97	N/A	N/A	75-12	0	N/A	0-20	
1,2,4-Trichlorobenzene	50.00	49.14	98	N/A	N/A	65-13	5	N/A	0-20	
1,1,1-Trichloroethane	50.00	46.95	94	N/A	N/A	65-13	0	N/A	0-20	
Hexachloro-1,3-Butadiene	50.00	46.63	93	N/A	N/A	50-14	0	N/A	0-20	



# **Quality Control - LCS/LCSD**

Environmental Science International, Inc.

Date Received:

Work Order:

14-07-1550

Kailua, HI 96734-2500

Preparation:

Method:

O7/23/14

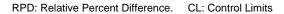
Work Order:

14-07-1550

GC/MS / EPA 8260B

Project: Red Hill LTM 112066 Page 5 of 5

<u>Parameter</u>	Spike Added	LCS Conc.	<u>LCS</u> %Rec.	LCSD Conc.	LCSD %Rec.	%Rec. CL	RPD	RPD CL	<u>Qualifiers</u>
1,1,2-Trichloroethane	50.00	50.31	101	N/A	N/A	75-125	N/A	0-20	
Trichloroethene	50.00	49.97	100	N/A	N/A	70-125	N/A	0-20	
1,2,3-Trichloropropane	50.00	48.66	97	N/A	N/A	75-125	N/A	0-20	
Vinyl Chloride	50.00	45.04	90	N/A	N/A	50-145	N/A	0-20	
p/m-Xylene	100.0	98.26	98	N/A	N/A	75-130	N/A	0-20	
o-Xylene	50.00	51.86	104	N/A	N/A	80-120	N/A	0-20	
Methyl-t-Butyl Ether (MTBE)	50.00	47.49	95	N/A	N/A	65-125	N/A	0-20	
Gasoline Range Organics	1000	1017	102	1007	101	80-120	1	0-20	





# **Sample Analysis Summary Report**

Work Order: 14-07-1550				Page 1 of 1
Method	Extraction	Chemist ID	Instrument	Analytical Location
EPA 6020	EPA 3005A Filt.	598	ICP/MS 04	1
EPA 8015B (M)	EPA 3510C	628	GC 45	1
EPA 8270C SIM PAHs	EPA 3510C	923	GC/MS AAA	1
GC/MS / EPA 8260B	EPA 5030C	849	GC/MS OO	2

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841 Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841



## **Glossary of Terms and Qualifiers**

Work Order: 14-07-1550 Page 1 of 1

<ul> <li>Qualifiers Definition</li> <li>* See applicable analysis comment.</li> <li>&lt; Less than the indicated value.</li> <li>&gt; Greater than the indicated value.</li> </ul>	control due to a required sample dilution. Therefore, the sample data was reported without further control due to matrix interference. The associated method blank surrogate spike compound was
<ul> <li>Less than the indicated value.</li> </ul>	control due to matrix interference. The associated method blank surrogate spike compound was
	control due to matrix interference. The associated method blank surrogate spike compound was
> Greater than the indicated value	control due to matrix interference. The associated method blank surrogate spike compound was
ordator trial trib irraidatou vardo.	control due to matrix interference. The associated method blank surrogate spike compound was
<ol> <li>Surrogate compound recovery was out of clarification.</li> </ol>	control due to matrix interference. The associated method blank surrogate spike compound was
Surrogate compound recovery was out of in control and, therefore, the sample data	was reported without further clarification.
3 Recovery of the Matrix Spike (MS) or Matr associated LCS recovery was in control.	ix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The
4 The MS/MSD RPD was out of control due	to suspected matrix interference.
5 The PDS/PDSD or PES/PESD associated	with this batch of samples was out of control due to suspected matrix interference.
6 Surrogate recovery below the acceptance	limit.
7 Surrogate recovery above the acceptance	limit.
B Analyte was present in the associated me	hod blank.
BU Sample analyzed after holding time expire	d.
BV Sample received after holding time expired	d.
DL The Detection Limit (DL) is the smallest ar the 99% level of confidence.	halyte concentration that can be demonstrated to be different from zero or a blank concentration at
E Concentration exceeds the calibration range	ge.
ET Sample was extracted past end of recomm	nended max. holding time.
HD The chromatographic pattern was inconsis	stent with the profile of the reference fuel standard.
ICH Initial calibration verification recovery is ab	ove the control limit for this analyte.
ICJ Initial calibration verification recovery is be	low the control limit for this analyte.
IH Calibration verification recovery is above t	ne control limit for this analyte.
IJ Calibration verification recovery is below the	ne control limit for this analyte.
J Analyte was detected at a concentration b	elow the LOQ and above the DL. Reported value is estimated.
JA Analyte positively identified but quantitatio	n is an estimate.
LOD The Limit of Detection (LOD) is the smalle detected at 99% confidence level.	st amount or concentration of a substance that must be present in a sample in order to be
LOQ The Limit of Quantitation (LOQ) is the low precision and bias.	est concentration of a substance that produces a quantitative result within specified limits of
Q Spike recovery and RPD control limits do concentration by a factor of four or greater	not apply resulting from the parameter concentration in the sample exceeding the spike

- SG The sample extract was subjected to Silica Gel treatment prior to analysis.
- Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD).
- X % Recovery and/or RPD out-of-range.
- Z Analyte presence was not confirmed by second column or GC/MS analysis.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

Calscience Environmental Laboratories, Inc. 

7440 Lincoln Way, Garden Grove, CA 92841-1427 • (7' Other locations: Concord, San Luis Obispo, Houston, and For courier service / sample drop off informatic contact sales@calscience.com or call us.

WO#/LAB USE (	
(714) 895-5494	and Corpus Christi nation,

CHAIN OF CUSTODY RECORD

Date Page\_

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ORATC	ADDRESS: 354 WILLING St., Suite 304	CITY Pailud	TEL:808-201-0740 FORDING RESCIENCE COM AFER	TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"):	SAME DAY	COELT EDF	SPECIAL INSTRUCTIONS:	Regular (Full) and monitoring analyte ist				<u> </u>	7.77	M	15	1,4-1		-	 	Relinquished by: (Signature)	Relinquished by: (Signature)	Relinquished by: (Signature)	oppusoodategoodoga	
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DISTRIBUTION: White with final report, Green and Yellow to Client.

Please note that pages 1 and 2 of 2 of our T/Cs are printed on the reverse side of the Green and Yellow copies respectively.

# Richard Villafania

From: Sent:

Ann Dang [ADang@esciencei.com] Tuesday, July 22, 2014 6:34 PM Richard Villafania

To:

Cc:

Domonkos Feher; Traci Sylva; Jeff Hattemer

Subject:

112066 Red Hill shipment 7-22-14

Attachments:

COC 7-22-14.pdf

Hi Richard,

We shipped 2 coolers today, the FedEx tracking is 804557917397 and 780047104871.

For ES107UF, "field filtered" should not be marked. The correction was made on the attached COC.

Thanks, Ann

Click here to report this email as spam.

01/01/14 Revision

130 140 CHAIN OF CUSTODY RECORD F 1200 SAMPLER(S): (PRINT) (3'002 N89 Time: Time: Time: 1712) 8.812 [] 6817 [] 3617 [] (IV)10 REQUESTED ANALYSES 12114 X747\0209 \Big X747\0109 \Big sisteM SST P.O. NO.: MIC 07S8 TO 0YS8 CD 2HA9 Please check box or fill in blank as needed. Date: Date: PCBs (8082) Pesticides (8081) Page\_ Date 800cs (8270) Prep (5035) 🗆 En Core 🗆 Terra Core Oxygenates (8260) **つつ021** CLIENT PROJECT NAME / NUMBER: ( hong (85e0) □ 81EX / MTBE □ 8260 □ WO # / LAB USE ONLY PROJECT CONTACT Hdl 24 1111 Poperty TPH C6-C36 C6-C44 Received by: (Signature/Affiliation) Received by: (Signature/Affiliation) Received by: (Signature/Affiliation (5/28 DECTED (b)Hard (1928) <del>- ОЯЭ-Б-</del> (в)нчт**Ж** TELSO8-2 COL. UTUD Churu PESCIEMET-CUM AFCHERBETTICKED.CAM WE Calscience Environmental Laboratories, Inc. Field Filtered CHES. CLETZY LOG CODE X Preserved × Χ LABORATORY CLIENT FAVI PO NYMEPIAI SCICNCE INTORNATIONA Other locations: Concord, San Luis Obispo, Houston, and Corpus Christi Unpreserved 7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494 X STANDARD S o S P S SNT. 9 σ For courier service / sample drop off information, contact sales@calscience.com or call us. REGULD (FULL) (AM MONITORING ANOTYTE 115) STATE: \*. 4 × 3 × 4 1. A CHOSE MATRIX 4 TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"): ☐5 DAYS 080 120 のなる 2501 MILETIC TIME 200 SAMPLING シミシ HIZZIL アルンバス ☐72 HR がんだし DATE 7 ☐48 HR ADDRESS: 3574 (ILMINU GLOBAL ID 5 107 UT SAMPLE ID Relinquished by: (Signature) ☐24 HR Relinquished by: (Signature) Relinquished by: (Signature) 557RM SPECIAL INSTRUCTIONS: んさせ ハヘころ 2106 OITY: KALTURA ☐ COELT EDF SAME DAY LAB USE ONLY

DISTRIBUTION: White with final report, Green and Yellow to Client. Please note that pages 1 and 2 of 2 of our T/Cs are printed on the reverse side of the Green and Yellow copies, respectively.





ORIGIN ID:HNLA (808) 261-0740 ESI

354 ULUNIU ST STE 304

KAILUA: HI 967342532 UNITED STATES US

SHIP DATE: 22JUL14 ACTWGT: 48.0 LB CAD: /POS1501 DIMS: 24x13x13 IN

TO SAMPLE CONTROL CAL SCIENCE 7440 LINCOLN WAY

# **GARDEN GROVE CA 92841**

1 of 2 TRK# 0200 8045 5791 7397

WED - 23 JUL AA



354 ULUNIU ST STE 304

KAILUA, HI 967342532 UNITED STATES US

BILL SENCER

**TO SAMPLE CONTROL** CAL SCIENCE 7440 LINCOLN WAY

# GARDEN GROVE CA 92841 (714) 895-5494 REF:

STANDARD OVERNIGHT

WZ APVA

92841 CA-US SNA





Calscience

WORK ORDER #: 14-07- ☐ 55 5 0

# SAMPLE RECEIPT FORM

Cooler \_\_\_ of 2

CLIENT: Env. Science.	DATE:	07/23	/14_
TEMPERATURE: Thermometer ID: SC1 (Criteria: 0.0 °C – 6.0 °C, not froze	n except se	ediment/tiss	ue)
	Blank	☐ Samp	
☐ Sample(s) outside temperature criteria (PM/APM contacted by:)			
☐ Sample(s) outside temperature criteria but received on ice/chilled on same c	lav of samp	lina.	
☐ Received at ambient temperature, placed on ice for transport by Co			
		Checked	by: <u>15</u>
Ambient Temperature:   Air   Filter		Onconoa	~ 3 ·
CUSTODY SEALS INTACT:			, 7
© Cooler □ □ No (Not Intact) □ Not Present	□ N/A	Checked I	by: 15
☑ Sample □ □ □ No (Not Intact) □ Not Present		Checked b	oy: <u>826</u>
			31/0
SAMPLE CONDITION:	Yes	No	N/A
Chain-Of-Custody (COC) document(s) received with samples			
COC document(s) received complete			***************************************
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.			
☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished.			
Sampler's name indicated on COC			
Sample container label(s) consistent with COC			
	,		
Proper containers and sufficient volume for analyses requested			. П
Analyses received within holding time	<i>y</i>		
Aqueous samples received within 15-minute holding time		П	
□ pH □ Residual Chlorine □ Dissolved Sulfides □ Dissolved Oxygen	A.		
Proper preservation noted on COC or sample container	. 🔑		لسنا
Unpreserved vials received for Volatiles analysis  Volatile analysis container(s) free of headspace		П	
Tedlar bag(s) free of condensation  CONTAINER TYPE:			استغي
Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve () □EnCore	es <sup>®</sup> □Terra	aCores® 🗆	
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Air: □Tedlar <sup>®</sup> □Canister Other: □ Trip Blank Lot#: 131007B		d/Checked b	
Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: E  Preservative: h: HCL n: HNO <sub>3</sub> na <sub>2</sub> :Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> na: NaOH p: H <sub>3</sub> PO <sub>4</sub> s: H <sub>2</sub> SO <sub>4</sub> u: Ultra-pure znna: ZnAc <sub>2</sub> +Na	•	Reviewed by Scanned b	-nl



# Calscience

WORK ORDER #: 14-07- [ 5 5 5

# SAMPLE RECEIPT FORM

Cooler  $\underline{\mathcal{L}}$  of  $\underline{\mathbf{Z}}$ 

CLIENT: Env. Science. DATE:	07/23/	14
TEMPERATURE: Thermometer ID: SC1 (Criteria: 0.0 °C – 6.0 °C, not frozen except se  Temperature 2 °C - 0.3 °C (CF) = 2 • 5 °C Blank  Sample(s) outside temperature criteria (PM/APM contacted by:)  Sample(s) outside temperature criteria but received on ice/chilled on same day of sampli	☐ Sample	
☐ Received at ambient temperature, placed on ice for transport by Courier.		
Ambient Temperature: □ Air □ Filter	Checked by	: 15
CUSTODY SEALS INTACT:  ☐ Cooler ☐ ☐ No (Not Intact) ☐ Not Present ☐ N/A ☐ Sample ☐ ☐ No (Not Intact) ☐ Not Present ☐ N/A	Checked by:	
SAMPLE CONDITION: Yes	No	N/A
Chain-Of-Custody (COC) document(s) received with samples		
COC document(s) received complete		
☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels.		
$\square$ No analysis requested. $\square$ Not relinquished. $\square$ No date/time relinquished.		
Sampler's name indicated on COC		
Sample container label(s) consistent with COC		
Sample container(s) intact and good condition		
Proper containers and sufficient volume for analyses requested		
Analyses received within holding time		
Aqueous samples received within 15-minute holding time		
□ pH □ Residual Chlorine □ Dissolved Sulfides □ Dissolved Oxygen □		
Proper preservation noted on COC or sample container		
Unpreserved vials received for Volatiles analysis		
Volatile analysis container(s) free of headspace		
Tedlar bag(s) free of condensation		9
Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve () □EnCores® □Terrad	Cores <sup>®</sup> □	
Aqueous: VOA UVOAh UVOAna2 U125AGB U125AGBh U125AGBp 21AGB U		1AGRe
□500AGB □500AGJ □500AGJs □250AGB □250CGB □250CGBs □1PB [		
□250PB ☑250PBny□125PB □125PBznna □100PJ □100PJna₂ □ □		
Air: □Tedlar <sup>®</sup> □Canister Other: □ Trip Blank Lot#: Labeled/		82h
Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope R	Reviewed by: _	<del>7</del> 28
Preservative: h: HCL n: HNO3 na2:Na2S2O3 na: NaOH p: H3PO4 s: H2SO4 u: Ultra-pure znna: ZnAc2+NaOH f: Filtered	Scanned by:	778

# **RAW DATA SHEET** FOR METHOD: EPA 8015B (M)

WORK ORDER:

D/T EXTRACTED:

14-07-1550

GC 45

INSTRUMENT: **EXTRACTION:** 

**EPA 3510C** 2014-07-24 00:00 ANALYZED BY: D/T ANALYZED:

628

2014-07-25 09:09

**REVIEWED BY:** 

D/T REVIEWED:

DATA FILE:

W:\GC 45\GC 45 DATA\2014\140724\14072455.D\14072455

#

**CLIENT SAMPLE NUMBER: ES106** 

LCS/MB BATCH: MS/MSD BATCH:

ug/L

140724B11B

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

FINAL VOLUME / WEIGHT:

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

ADJUSTMENT RATIO TO PF:

0.50

COMMENT:

UNITS:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

É

8

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND TPH as Diesel

INI. CONC

7390

DF 1.00 CONC 37.0

4 5 <u>D\</u>

101 : A &

<u>DL</u> 11

LOD 12

LOQ **QUAL** 

25

b

\_\_\_\_\_\_

## Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072455.D

Vial Number : Vial 55

Page Number : 1
Operator : 682 Vial Number : Vial Instrument : GC 45 Injection Number : 1
Sample Name : 14-07-1550-1 Sequence Line : 55

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 25 Jul 14 9:09:52 AM

Report Created on: 25 Jul 14 06:12 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

Sig. 1 in W:\GC 45\GC 45 DATA\2014\140724\ ->

Pk	Ret Time	\GC 45 DATA\2014 Area	Height	Peak	Width	Response %
		0.27		   VV 0	0.025	0.029
2		0.27		0 VV	0.024	0.029
3		0.98		0 VV	0.046	0.107
4		1.57		1 VV	0.030	0.171
Ē		0.52		0 VV	0.024	0.056
6		0.50		0 VV	0.023	0.055
7		1.30	•	1 VV	0.040	0.141
8		2.38		1 VV	0.043	0.259
9		1.42		1 VV	0.041	0.155
10		1.97	• •	1 VV	0.034	0.214
13		1.57		1 VV	0.031	0.170
12		4.47		1 VV	0.050	0.486
13		6.75	Ţ.	3 VV	0.037	0.734
14		4.63		1 VV	0.053	0.504
1.5		4.30	Apr	2 VV	0.042	0.468
16	4.226	5.53		2 VV ,	0.051	0.602
1.	7 4.285	7.70		2 VV ′	0.061	0.837
18	3 4.327	1.85		2 VV	0.015	0.201
19	9 4.393	8.42	•	2 VV	0.062	0.916
20	4.413	2.78		2 VV	0.020	0.303
2:	1.4.430	10.16		3 VV	0.054	1.105
22	4.489	3.55		2 VV	0.026	0.386
23	3 4.511	5.51		2 VV	0.040	0.599
24	4.565	9.84		2 VV	0.072	1.070
2		4.25		2 VV	0.032	0.462
26	4.660	3.64	-	2 VV	0.028	0.396
2	7 4.726	23.63		6 VV	0.068	2.570
2		2.64		2 VV.	0.024	0.287
2		2.76		2 VV	0.025	0.301
3		4.16		2 VV	0.039	0.452
3		4.46		2 VV	0.044	0.485
3		1.55		1 VV	0.019	0.168
3		3.42		1 VV	0.041	0.372
3		2.97			0.038	0.323
3		5.93	- No	T 55	0.054	0.645
3		2.76		1 VVq	0.054	0.300
3		3.08		1 VV	0.038	
3		1.92	7.	1 VV	0.048	0.209
3	9 5.378	763.94		549 VV	0.023	83.096

Total area = 919.34

### Area Percent Report

: W:\GC\_45\GC 45 DATA\2014\140724\14072455.D Data File Name

: 2 Page Number

Sample Name

: 682 Operator Instrument : GC 45

: 14-07-1550-1

Vial Number : Vial 55

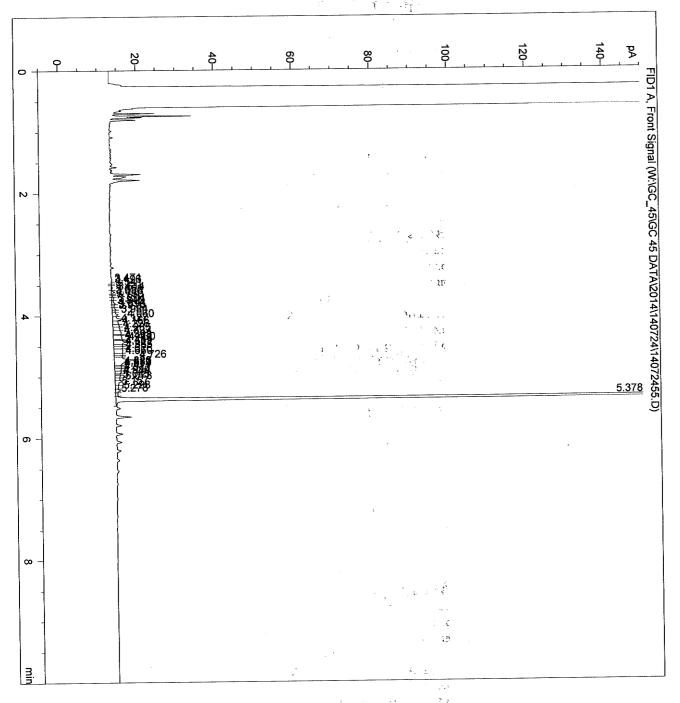
Injection Number: 1 Sequence Line : 55

Instrument Method: C:\CHEM32\1\METHODS\

9:09:52 AM : 25 Jul 14 Acquired on

Analysis Method : 8015B.MTH Report Created on: 25 Jul 14 06:12 pm

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies



# **RAW DATA SHEET** FOR METHOD: EPA 8015B (M)

**WORK ORDER: INSTRUMENT:** 

14-07-1550

GC 45

**EPA 3510C EXTRACTION:** 

D/T EXTRACTED: 2014-07-24 00:00 ANALYZED BY:

628 D/T ANALYZED:

2014-07-25 09:28

**REVIEWED BY:** 

D/T REVIEWED:

DATA FILE:

W:\GC 45\GC 45 DATA\2014\140724\14072456.D\14072456

2

**CLIENT SAMPLE NUMBER: ES107** 

LCS/MB BATCH:

140724B11B MS/MSD BATCH:

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

FINAL VOLUME / WEIGHT:

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

ug/L

ADJUSTMENT RATIO TO PF:

0.50

COMMENT:

UNITS:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND TPH as Diesel

INI. CONC

DF

CONC

<u>DL</u>

LOD 12

QUAL LOQ

107

1.00

ND

311

36.

11

25

Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072456.D

Page Number : 1

Operator : 682 Vial Number : Vial 56

Instrument : GC 45 Injection Number : 1
Sample Name : 14-07-1550-2 Sequence Line : 56

Instrument Method: C:\CHEM32\1\METHODS\

Acquired on : 25 Jul 14 9:28:40 AM

Report Created on: 25 Jul 14 06:12 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

Sig. 1 in W:\GC\_45\GC 45 DATA\2014\140724\ ->

Pk	Ret Time	Area				Response %
. 1	4.063	2.25	2	ΛŃ	0.019	0.290
2	5.394	771.63	549	VB	0.021	99.710

Total area = 773.87

#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072456.D

Page Number : 2

Operator : 682 Vial Number : Vial 56

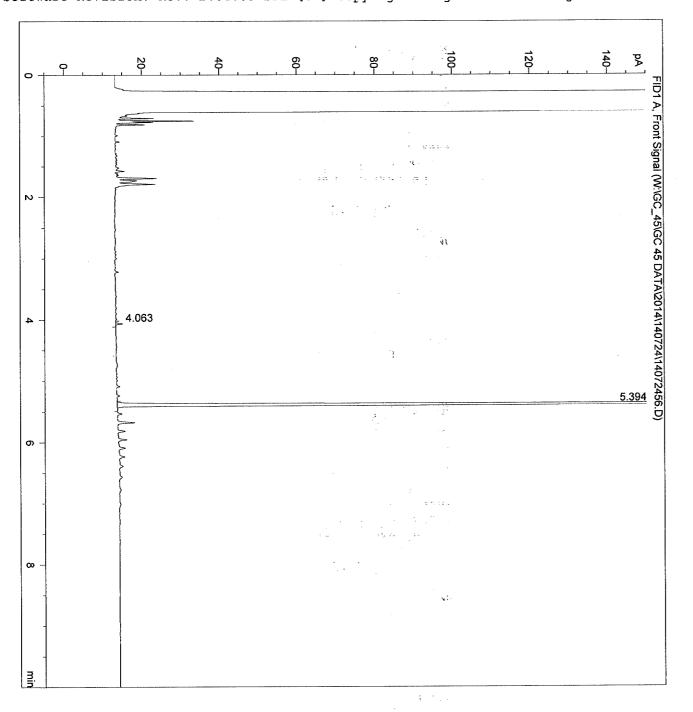
Instrument : GC 45 Injection Number : 1 Sample Name : 14-07-1550-2 Sequence Line : 56

Instrument Method: C:\CHEM32\1\METHODS\ -:

Acquired on : 25 Jul 14 9:28:40 AM

Report Created on: 25 Jul 14 06:12 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies



# **RAW DATA SHEET** FOR METHOD: EPA 8015B (M)

**WORK ORDER:** 

14-07-1550

**INSTRUMENT:** 

GC 45

**EXTRACTION:** 

**EPA 3510C** 

140724B11B

D/T EXTRACTED: 2014-07-24 00:00 ANALYZED BY:

D/T ANALYZED:

628 2014-07-25 09:47

**REVIEWED BY:** D/T REVIEWED:

DATA FILE:

W:\GC 45\GC 45 DATA\2014\140724\14072457.D\14072457

**CLIENT SAMPLE NUMBER: ES108** 

LCS/MB BATCH:

MS/MSD BATCH: UNITS:

ug/L

SAMPLE VOLUME / WEIGHT:

FINAL VOLUME / WEIGHT: ADJUSTMENT RATIO TO PF: DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

0.50

**COMMENT:** 

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND

TPH as Diesel

INI. CONC 270

DF 1.00 CONC ND

DL 11

LOD 12

LOQ

QUAL

25

\_\_\_\_\_\_\_

#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072457.D

Page Number : 1

Operator : 682 Vial Number : Vial 57

Instrument : GC 45 Injection Number : 1
Sample Name : 14-07-1550-3 Sequence Line : 57

Instrument Method: C:\CHEM32\1\METHODS\ ->

.19 Y

THE

Acquired on : 25 Jul 14 9:47:30 AM

Report Created on: 25 Jul 14 06:12 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies

Sig. 1 in W:\GC\_45\GC 45 DATA\2014\140724\ ->

Pk	Ret Time	Area	Height			Response	
						-	
. 1	4.066	2.43	2	VB	0.020	0.28	9
2	4.741	2.28	1	VV	0.031	0.27	1
3	4.783	0.97	1	VV	0.028	0.11	5
4	5.408	837.46	593	VV	0.022	99.32	5

Total area = 843.14

#### Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072457.D

Page Number : 2

Operator : 682

Vial Number : Vial 57

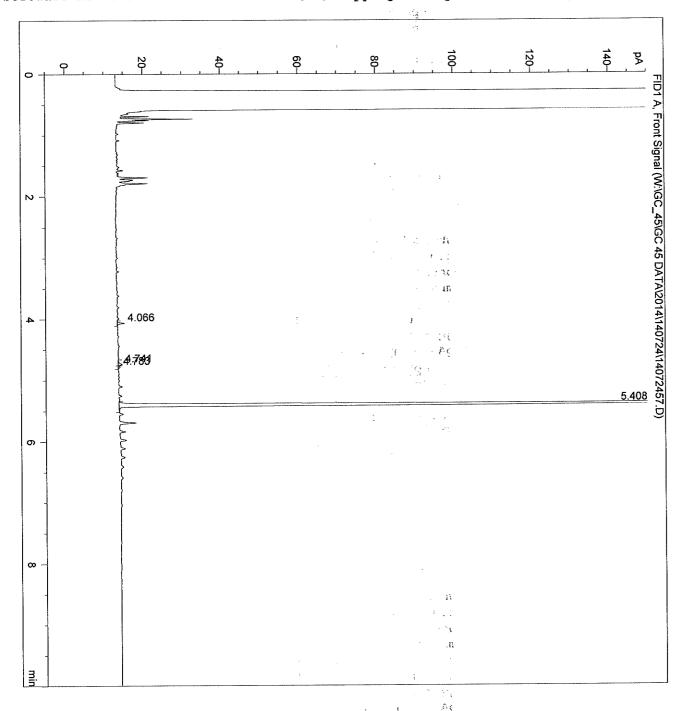
Instrument : GC 45 Injection Number : 1
Sample Name : 14-07-1550-3 Sequence Line : 57

Instrument Method: C:\CHEM32\1\METHODS\ -

Acquired on : 25 Jul 14 9:47:30 AM

Report Created on: 25 Jul 14 06:12 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies



# Area Percent Report

Data File Name : W:\GC\_45\GC 45 DATA\2014\140724\14072436.D

: 2 ?age Number : 682 )perator

Vial Number : Vial 36 Injection Number: 1

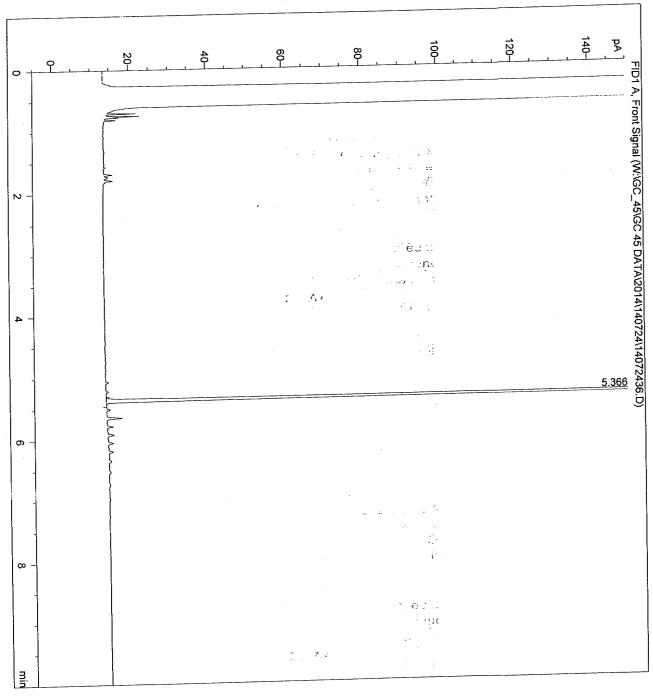
: GC 45 [nstrument : MB 14072411/12 Bample Name

Sequence Line : 36 Instrument Method: C:\CHEM32\1\METHODS\

: 25 Jul 14 3:30:24 AM Acquired on

Analysis Method : 8015B.MTH Report Created on: 25 Jul 14 06:11 pm

Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies



: W:\GC\_45\GC 45 DATA\2014\140724\14072434.D ata File Name

age Number : 682 )perator

: Vial 34 Vial Number

Injection Number : 1

Instrument Sample Name : GC 45 Sequence Line : D400 C28 50 L041814D

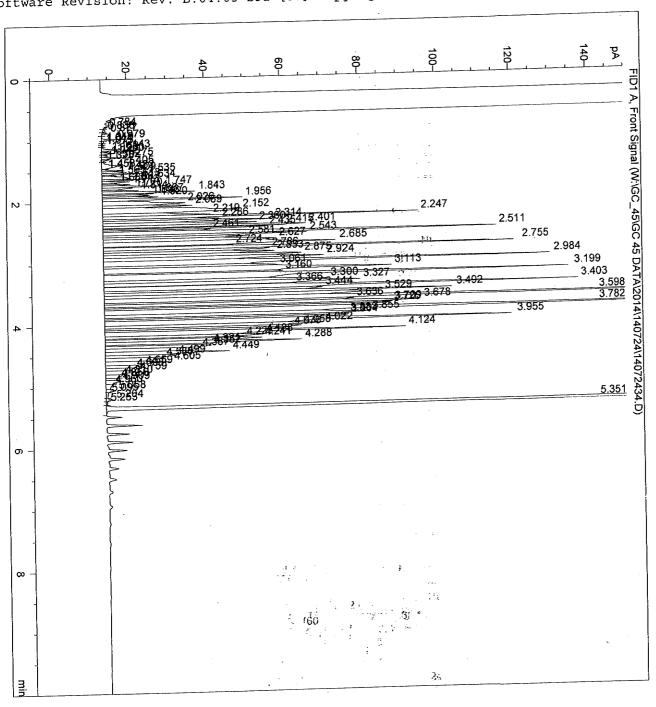
Instrument Method: C:\CHEM32\1\METHODS\

cquired on

2:53:13 AM : 25 Jul 14

leport Created on: 25 Jul 14 06:11 pm Analysis Method : 8015B.MTH

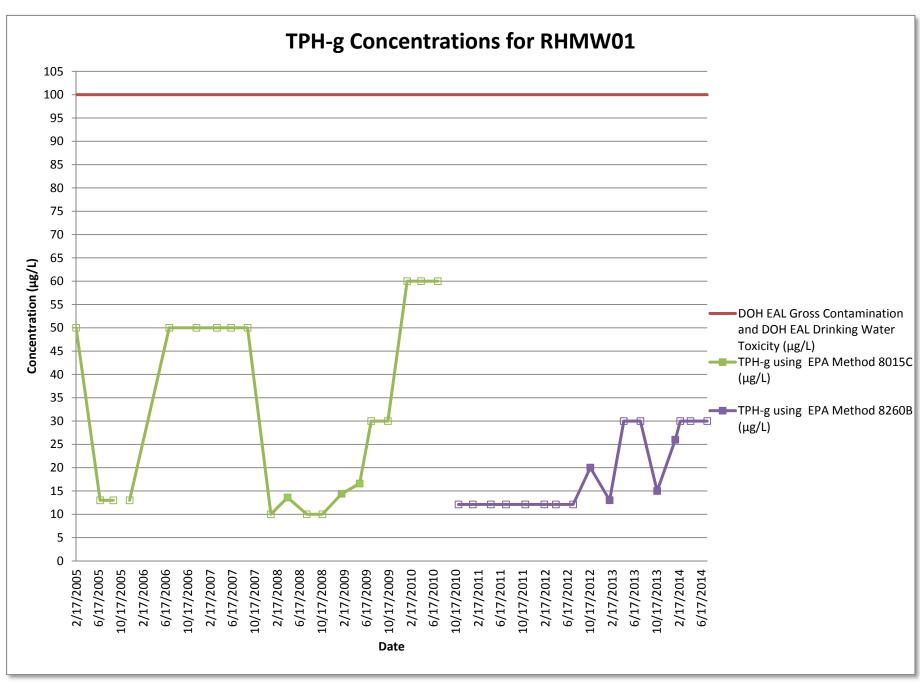
Software Revision: Rev. B.04.03-SP1 [87] Copyright © Agilent Technologies





## APPENDIX D Historical Groundwater Exceedance Trends

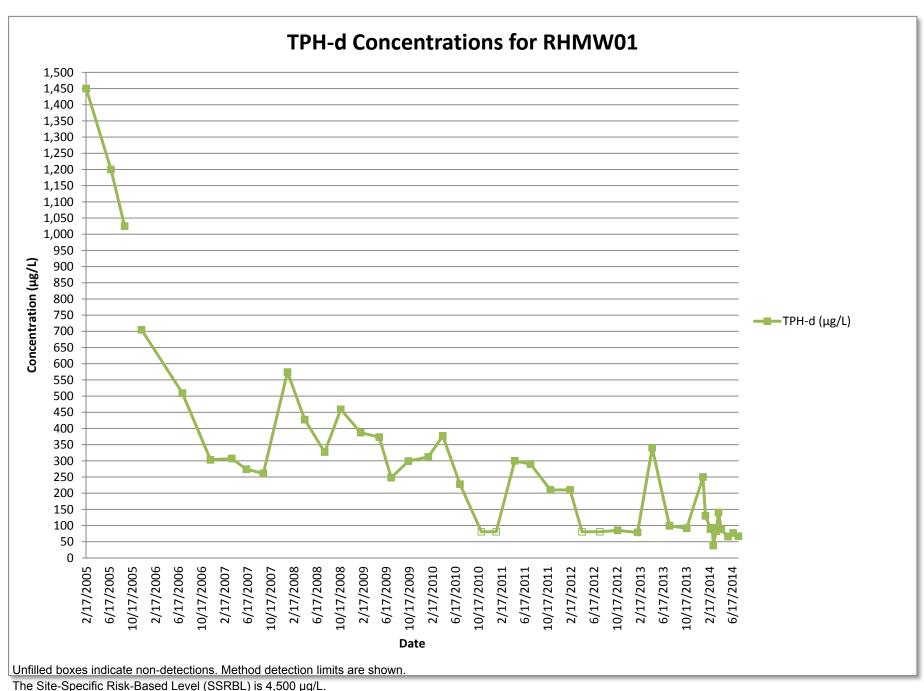




Data points for 2/17/2005 through 9/8/2005 and 12/6/2005 are the average of the primary and duplicate samples. Possible laboratory contamination for 10/22/2012, 10/21/2013, and 1/28/2014 sampling events.

Unfilled boxes indicate non-detections. Method detection limits are shown.

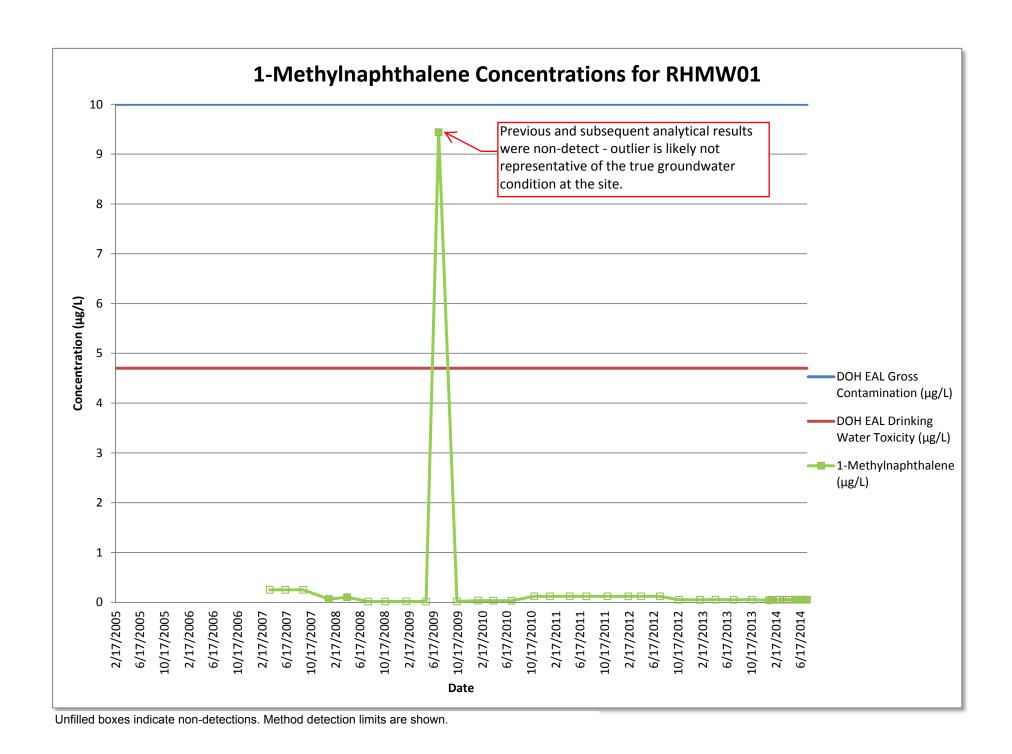




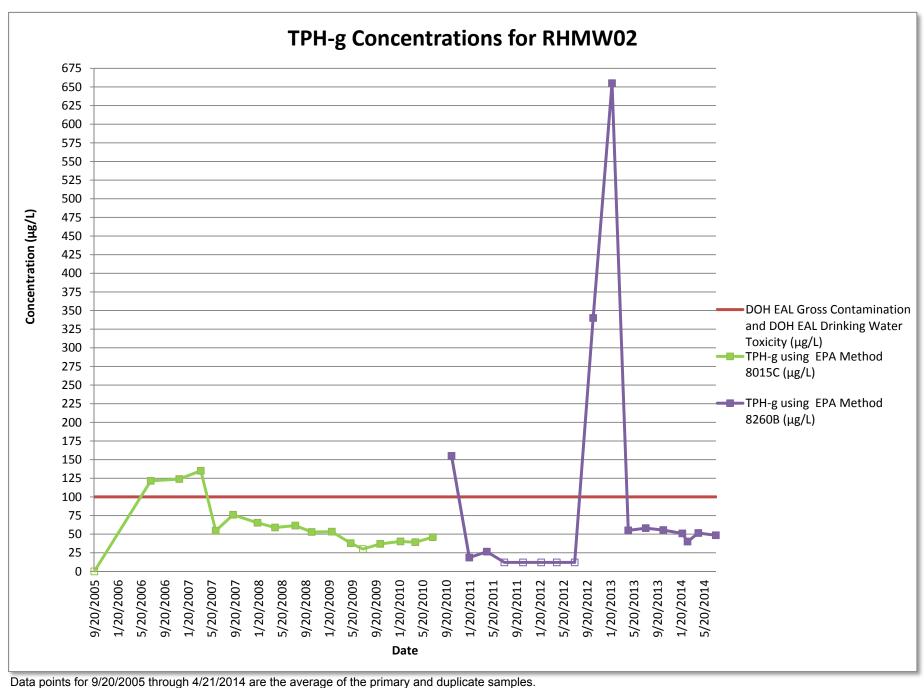
Numerous sample results had a chromatographic pattern that didn't match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

Data points for 2/17/2005 through 9/8/2005 and 12/6/2005 are the average of the primary and duplicate samples.







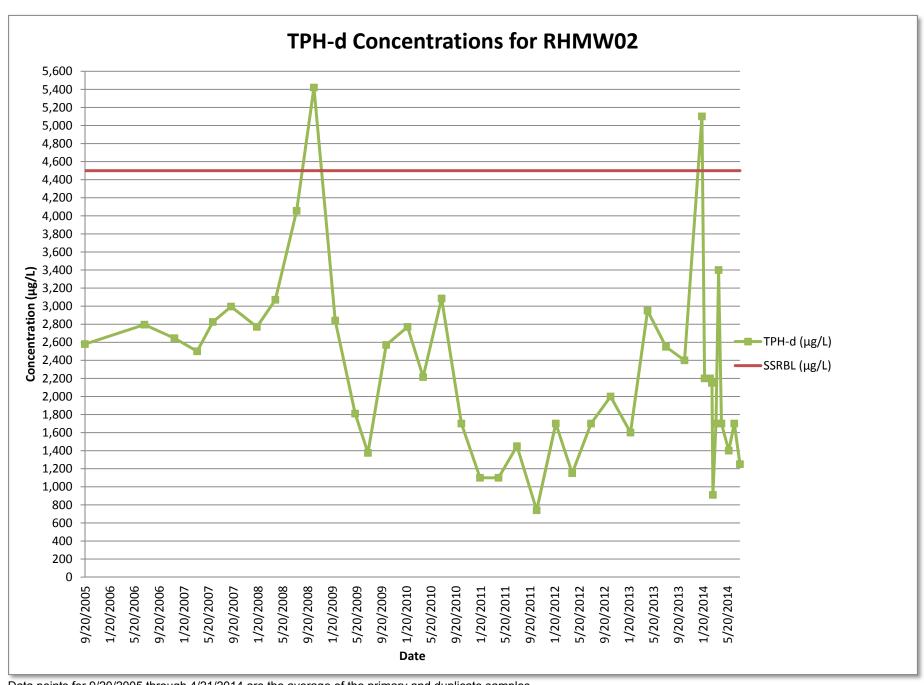


Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events.

Unfilled boxes indicate non-detections. Method detection limits are shown.

Primary sample results are shown for 1/26/2012 and 7/18/2012; all other concentrations are the average of the primary and duplicate sample results.



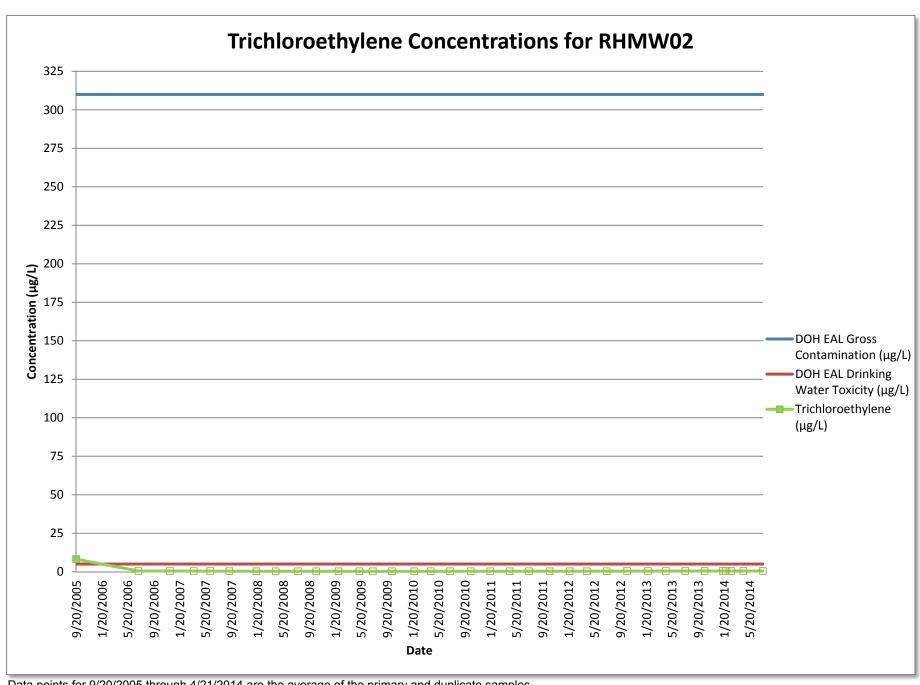


Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples.

Unfilled boxes indicate non-detections. Method detection limits are shown.

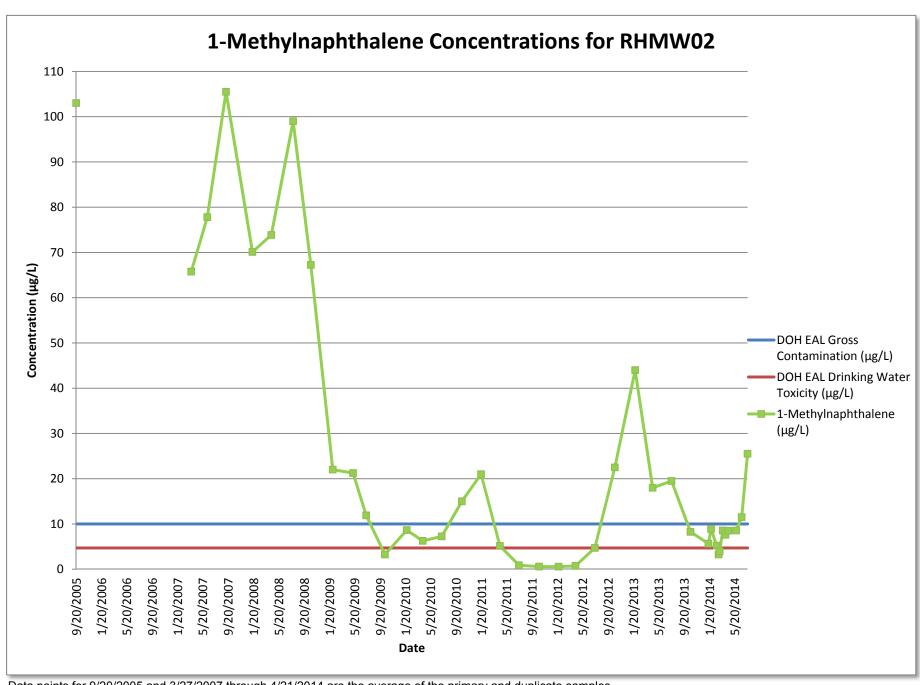
Numerous sample results had a chromatographic pattern that didn't match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.





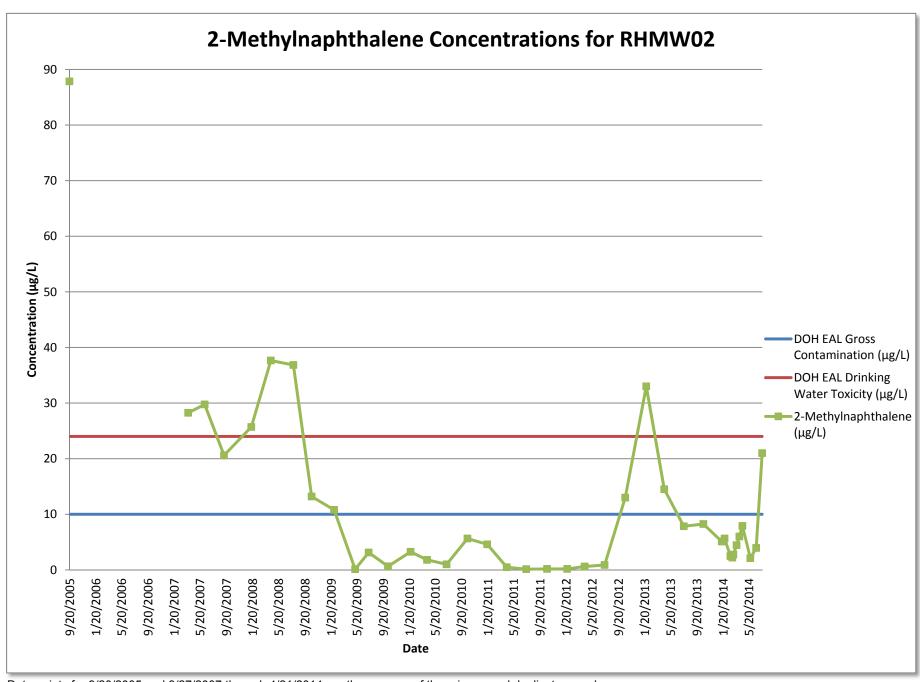
Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples. Unfilled boxes indicate non-detections. Method detection limits are shown.





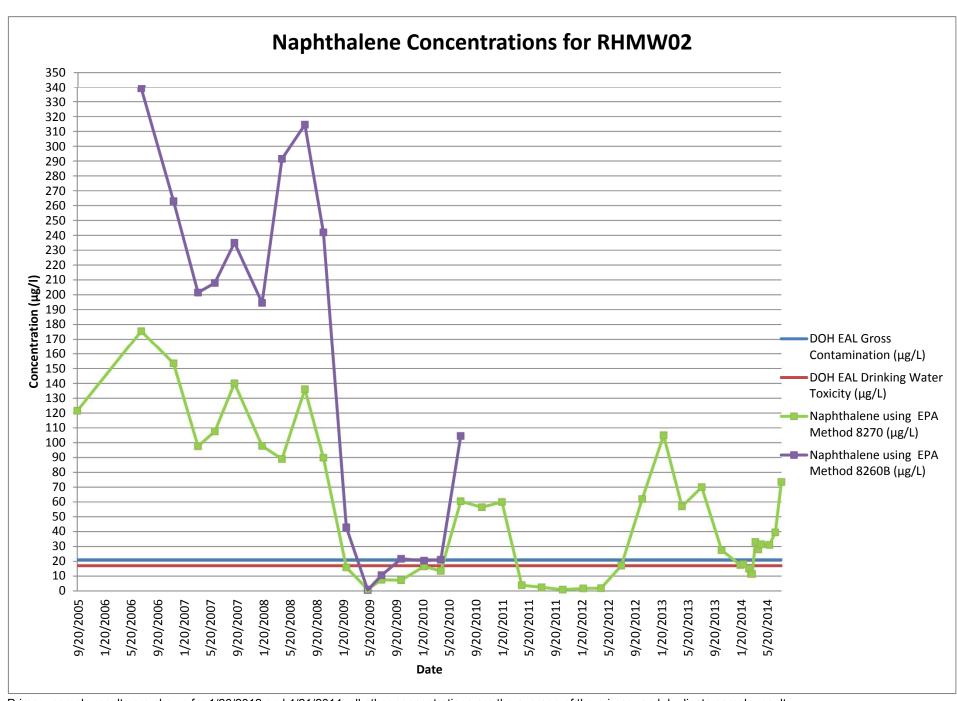
Data points for 9/20/2005 and 3/27/2007 through 4/21/2014 are the average of the primary and duplicate samples. Unfilled boxes indicate non-detections. Method detection limits are shown.





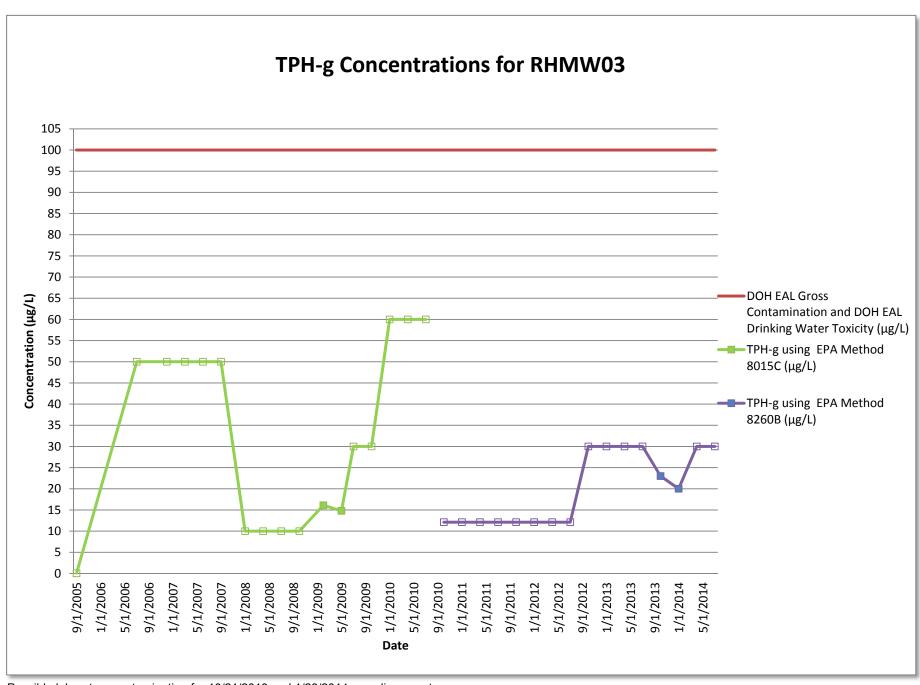
Data points for 9/20/2005 and 3/27/2007 through 4/21/2014 are the average of the primary and duplicate samples. Unfilled boxes indicate non-detections. Method detection limits are shown.





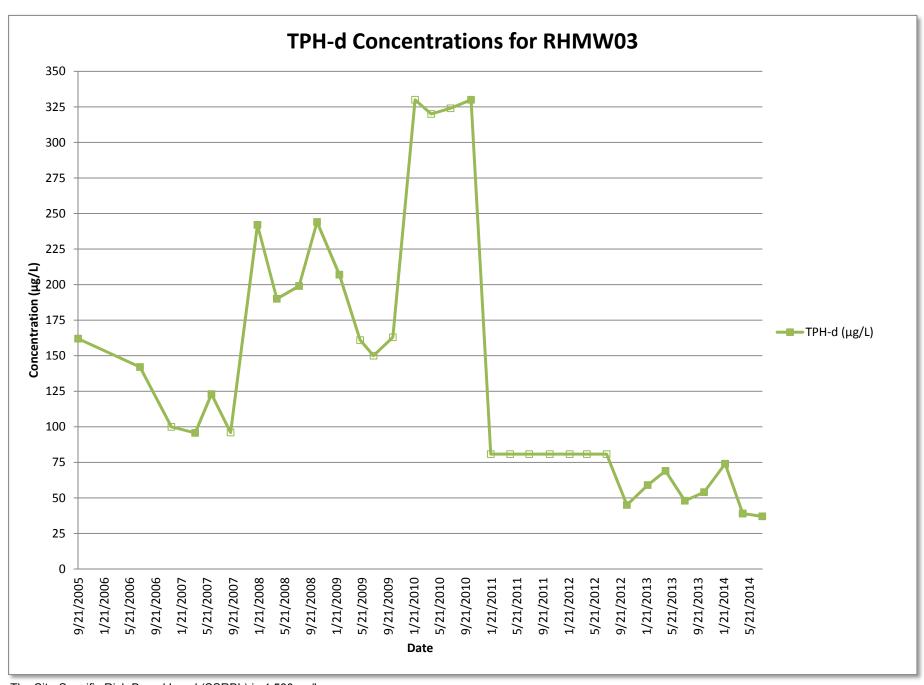
Primary sample results are shown for 1/26/2012 and 4/21/2014; all other concentrations are the average of the primary and duplicate sample results. Unfilled boxes indicate non-detections. Method detection limits are shown.





Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events. Unfilled boxes indicate non-detections. Method detection limits are shown.



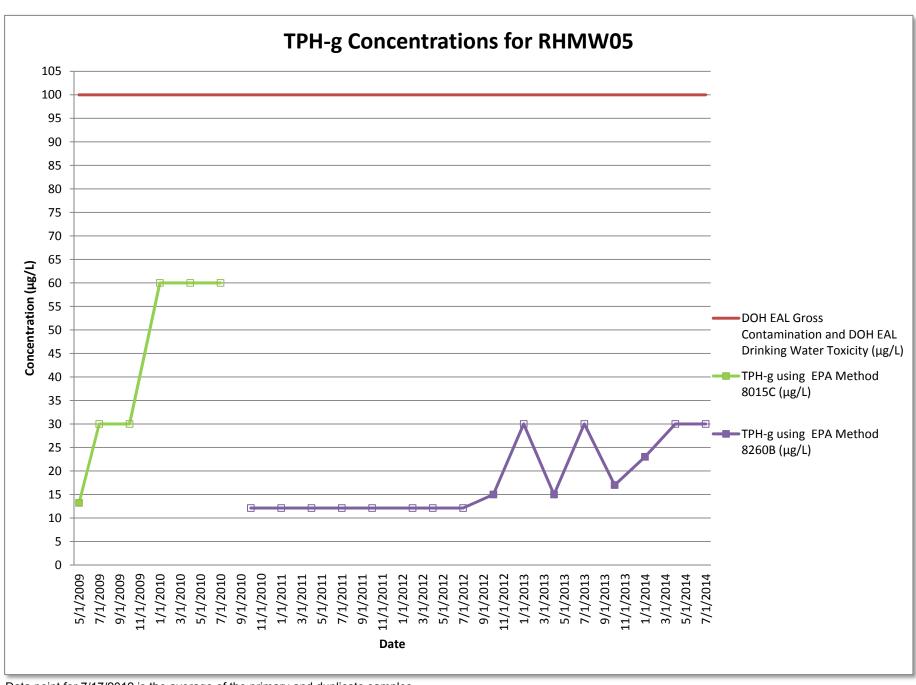


The Site-Specific Risk-Based Level (SSRBL) is 4,500 µg/L.

Unfilled boxes indicate non-detections. Method detection limits are shown.

Numerous sample results had a chromatographic pattern that didn't match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.



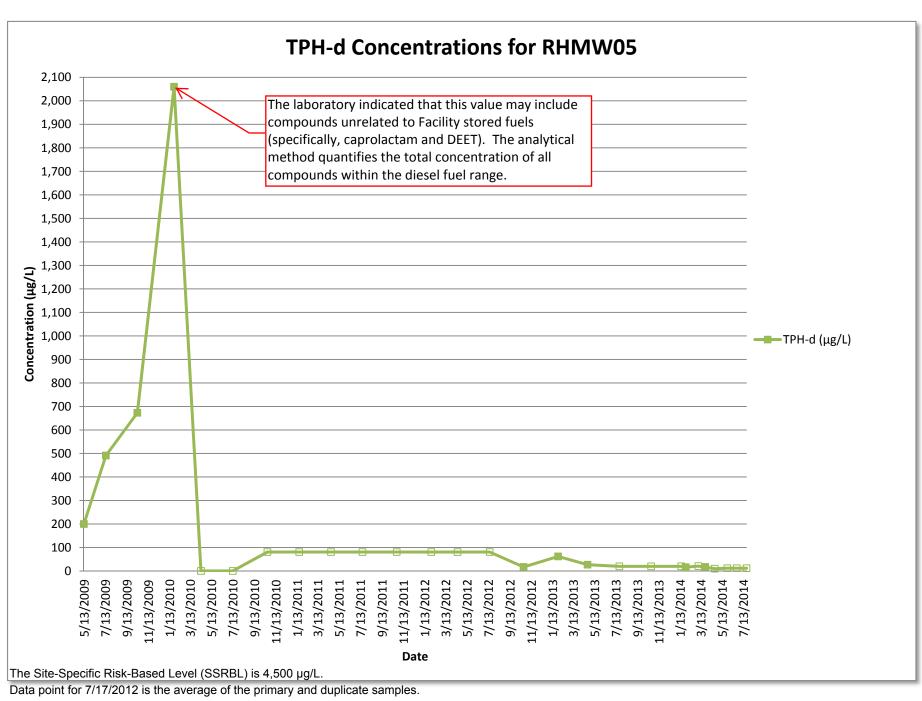


Data point for 7/17/2012 is the average of the primary and duplicate samples.

Unfilled boxes indicate non-detections. Method detection limits are shown.

Possible laboratory contamination for 10/22/2012, 10/22/2013, and 1/29/2014 sampling events.

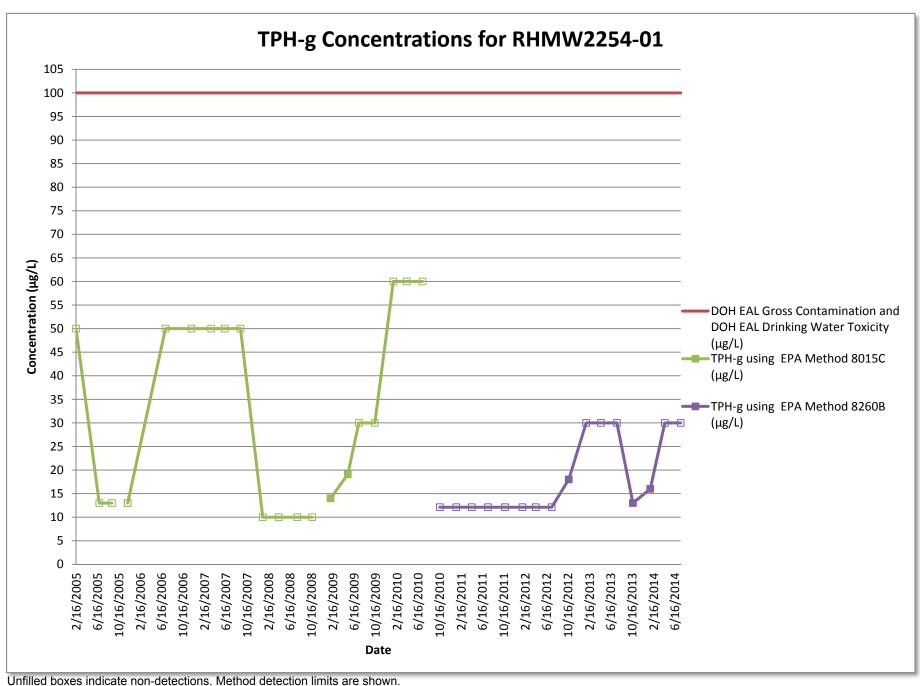




Unfilled boxes indicate non-detections. Method detection limits are shown.

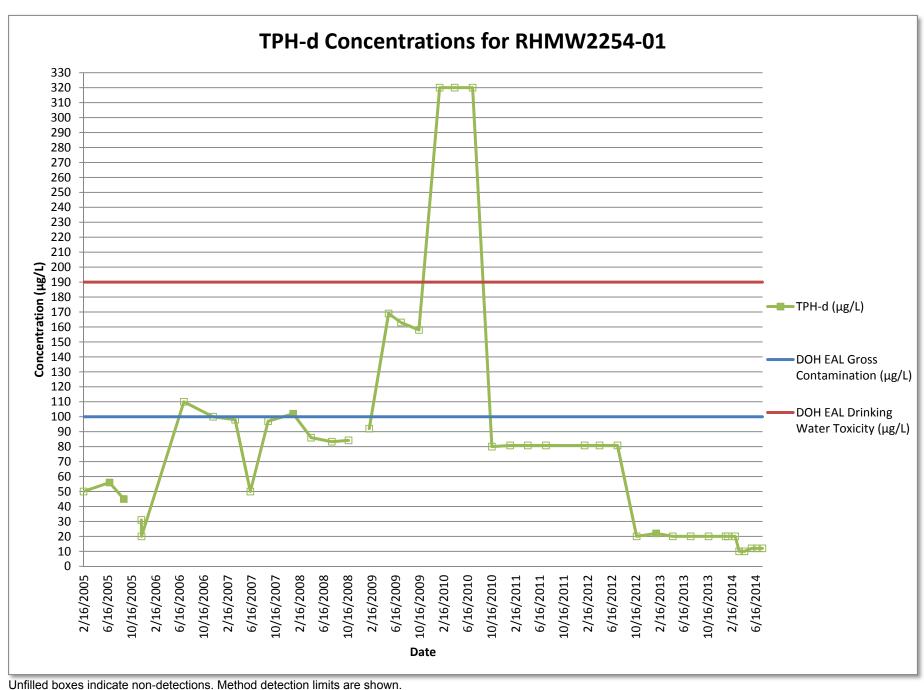
Numerous sample results had a chromatographic pattern that didn't match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.





Possible laboratory contamination for 10/22/2012, 10/22/2013, and 1/29/2014 sampling events.





Laboratory data rejected for 1/15/2008 sampling event.

Numerous sample results had a chromatographic pattern that didn't match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

